

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
 Data File : BN033865.D
 Acq On : 11 Sep 2024 17:35
 Operator : JU/RC
 Sample : P3878-18
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DDCC1

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BN033865.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.569	432	439	447	rVB2	2787511	4487339	1.96%	0.323%
2	6.581	606	611	614	rBV	3105737	4723883	2.07%	0.340%
3	6.687	621	629	635	rBV4	3819794	8365001	3.66%	0.602%
4	6.957	670	675	680	rBV	4250391	6807542	2.98%	0.490%
5	7.157	702	709	717	rBV2	16232489	28245355	12.35%	2.034%
6	7.498	761	767	771	rVB2	4778448	7768308	3.40%	0.559%
7	7.587	777	782	785	rBV	4219027	6817292	2.98%	0.491%
8	7.775	810	814	824	rVB6	2345212	6469148	2.83%	0.466%
9	7.934	836	841	846	rBV	5598149	8760365	3.83%	0.631%
10	7.993	846	851	855	rBV	7104887	10376733	4.54%	0.747%
11	8.046	855	860	865	rVB2	6031153	9804529	4.29%	0.706%
12	8.104	865	870	873	rBV	3458033	5353228	2.34%	0.386%
13	8.140	873	876	879	rBV2	3968098	5665164	2.48%	0.408%
14	8.175	879	882	889	rVB	4950453	7843934	3.43%	0.565%
15	8.275	889	899	901	rBV2	4348592	7478607	3.27%	0.539%
16	8.298	901	903	910	rVB	4281308	6595083	2.88%	0.475%
17	8.451	924	929	932	rBV	2796943	4004298	1.75%	0.288%
18	8.498	932	937	944	rVB	9237551	14165526	6.20%	1.020%
19	8.604	945	955	956	rBV	4361382	7887922	3.45%	0.568%
20	8.740	972	978	984	rVB	13655558	22506510	9.84%	1.621%
21	8.845	987	996	1003	rVB7	2843901	7312675	3.20%	0.527%
22	8.928	1003	1010	1014	rBV2	2078041	4074151	1.78%	0.293%
23	9.357	1078	1083	1090	rBV3	1747387	4250126	1.86%	0.306%
24	9.575	1110	1120	1126	rVB4	1490937	4308301	1.88%	0.310%
25	10.322	1230	1247	1251	rBV2	27093345	85780229	37.52%	6.177%
26	10.422	1255	1264	1269	rBV	18061536	31952310	13.97%	2.301%
27	10.569	1284	1289	1294	rBV2	3707095	6534088	2.86%	0.471%
28	10.669	1300	1306	1314	rVB2	3801386	7839916	3.43%	0.565%
29	10.804	1323	1329	1335	rVV	13377164	23270843	10.18%	1.676%
30	11.122	1376	1383	1387	rBV4	2021183	4589168	2.01%	0.330%
31	11.175	1387	1392	1395	rVV	6094712	10493683	4.59%	0.756%
32	11.204	1395	1397	1405	rVB2	4548072	5954775	2.60%	0.429%
33	11.322	1411	1417	1422	rBV3	7349282	11601883	5.07%	0.836%
34	11.381	1422	1427	1431	rBV	5956325	9571842	4.19%	0.689%
35	11.528	1446	1452	1454	rBV	4723732	7866685	3.44%	0.567%
36	11.557	1454	1457	1462	rVV	6780805	9650977	4.22%	0.695%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
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37	11.639	1462	1471	1476	rVB	3565806	6627819	2.90%	0.477%
38	11.710	1476	1483	1488	rBV4	7849586	16502333	7.22%	1.188%
39	11.757	1488	1491	1495	rVV	4591228	6266315	2.74%	0.451%
40	11.939	1518	1522	1531	rVB	4232499	7064482	3.09%	0.509%
41	12.145	1550	1557	1566	rVB2	19847956	37386093	16.35%	2.692%
42	12.392	1594	1599	1608	rVV4	3917477	8983916	3.93%	0.647%
43	12.486	1608	1615	1620	rVV4	2643084	5775652	2.53%	0.416%
44	12.581	1621	1631	1638	rVV4	8252855	16172947	7.07%	1.165%
45	12.728	1648	1656	1661	rVB2	8070561	14783506	6.47%	1.065%
46	12.992	1694	1701	1705	rVV2	7517927	13795690	6.03%	0.993%
47	13.039	1705	1709	1712	rVV	2644855	4035202	1.76%	0.291%
48	13.210	1733	1738	1746	rVB3	8424551	15313377	6.70%	1.103%
49	13.304	1747	1754	1756	rBV3	3930246	6534763	2.86%	0.471%
50	13.469	1776	1782	1786	rVV	4343784	8363062	3.66%	0.602%
51	13.516	1786	1790	1795	rVV4	6055736	10685934	4.67%	0.770%
52	13.569	1795	1799	1804	rVB	3714833	5195462	2.27%	0.374%
53	13.657	1807	1814	1818	rVV3	2393685	4482246	1.96%	0.323%
54	13.704	1818	1822	1826	rVV3	3768085	5166417	2.26%	0.372%
55	13.798	1831	1838	1841	rBV4	4953084	10337980	4.52%	0.744%
56	13.904	1852	1856	1861	rVB	5970676	7816274	3.42%	0.563%
57	13.986	1865	1870	1875	rVV3	2528454	4438227	1.94%	0.320%
58	14.086	1881	1887	1890	rBV	14668862	21052089	9.21%	1.516%
59	14.251	1910	1915	1919	rBV2	4651803	6363904	2.78%	0.458%
60	14.310	1919	1925	1929	rBV	12109872	16566091	7.25%	1.193%
61	14.386	1929	1938	1942	rBV5	2829448	7692877	3.36%	0.554%
62	14.639	1976	1981	1985	rVB2	3068282	4700939	2.06%	0.339%
63	14.704	1986	1992	1996	rBV2	5295619	9501651	4.16%	0.684%
64	14.745	1996	1999	2001	rVV3	3669439	5174252	2.26%	0.373%
65	14.810	2003	2010	2018	rVB	24000586	52965782	23.16%	3.814%
66	14.933	2023	2031	2035	rBV2	11352180	22868417	10.00%	1.647%
67	15.045	2042	2050	2054	rVB2	12584625	21025109	9.20%	1.514%
68	15.151	2063	2068	2072	rVB2	4519161	8134512	3.56%	0.586%
69	15.280	2083	2090	2097	rBV2	9288948	17205416	7.52%	1.239%
70	15.445	2111	2118	2128	rBV2	7708099	17757377	7.77%	1.279%
71	15.598	2140	2144	2148	rBV2	2964772	4707284	2.06%	0.339%
72	15.804	2174	2179	2186	rBV5	5104735	9327279	4.08%	0.672%
73	15.910	2192	2197	2200	rBV	12977290	20305553	8.88%	1.462%
74	16.622	2303	2318	2329	rBV6	33518505	228652757	100.00%	16.466%
75	16.757	2337	2341	2350	rVB2	8689862	15744837	6.89%	1.134%
76	16.833	2351	2354	2362	rVB4	3046058	4020845	1.76%	0.290%

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Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF Filtering: 5
 Sampling : 1 Min Area: 0.1 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
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77	16.945	2368	2373	2377	rBV2	5620260	11121072	4.86%	0.801%
78	17.033	2384	2388	2392	rVB2	6415297	9123790	3.99%	0.657%
79	17.080	2392	2396	2400	rBV3	8555146	12237763	5.35%	0.881%
80	17.127	2400	2404	2407	rVV	4151635	5173184	2.26%	0.373%
81	17.245	2416	2424	2428	rVB5	6581654	13387681	5.86%	0.964%
82	17.445	2453	2458	2462	rVV	13412002	17668445	7.73%	1.272%
83	17.498	2465	2467	2475	rVB2	3890667	5523824	2.42%	0.398%
84	17.863	2522	2529	2539	rVB2	9355769	20101959	8.79%	1.448%
85	18.133	2571	2575	2580	rVB	11563272	14994317	6.56%	1.080%
86	18.774	2680	2684	2688	rVB	10276503	12580713	5.50%	0.906%
87	19.139	2742	2746	2755	rVB6	4340478	7505785	3.28%	0.541%
88	19.316	2771	2776	2780	rBV	5399775	7799636	3.41%	0.562%
89	19.363	2780	2784	2788	rVB	7777567	9847027	4.31%	0.709%
90	19.898	2871	2875	2880	rBV	10713662	12377835	5.41%	0.891%
91	20.398	2956	2960	2964	rBV	8379550	9822842	4.30%	0.707%
92	20.874	3036	3041	3050	rVB2	16817051	25930643	11.34%	1.867%
93	21.357	3119	3123	3130	rVB	7757463	10499244	4.59%	0.756%
94	21.780	3191	3195	3200	rVB	3820807	5440122	2.38%	0.392%
95	21.892	3208	3214	3221	rBV3	9188242	17176996	7.51%	1.237%
96	22.492	3311	3316	3323	rVB	4439859	7118746	3.11%	0.513%
97	23.180	3428	3433	3441	rBV2	3049066	6077166	2.66%	0.438%
98	23.739	3523	3528	3539	rVB	3320308	7837218	3.43%	0.564%
99	23.927	3554	3560	3565	rBV	2829033	6149166	2.69%	0.443%
100	23.986	3566	3570	3579	rVB	2104141	4429155	1.94%	0.319%

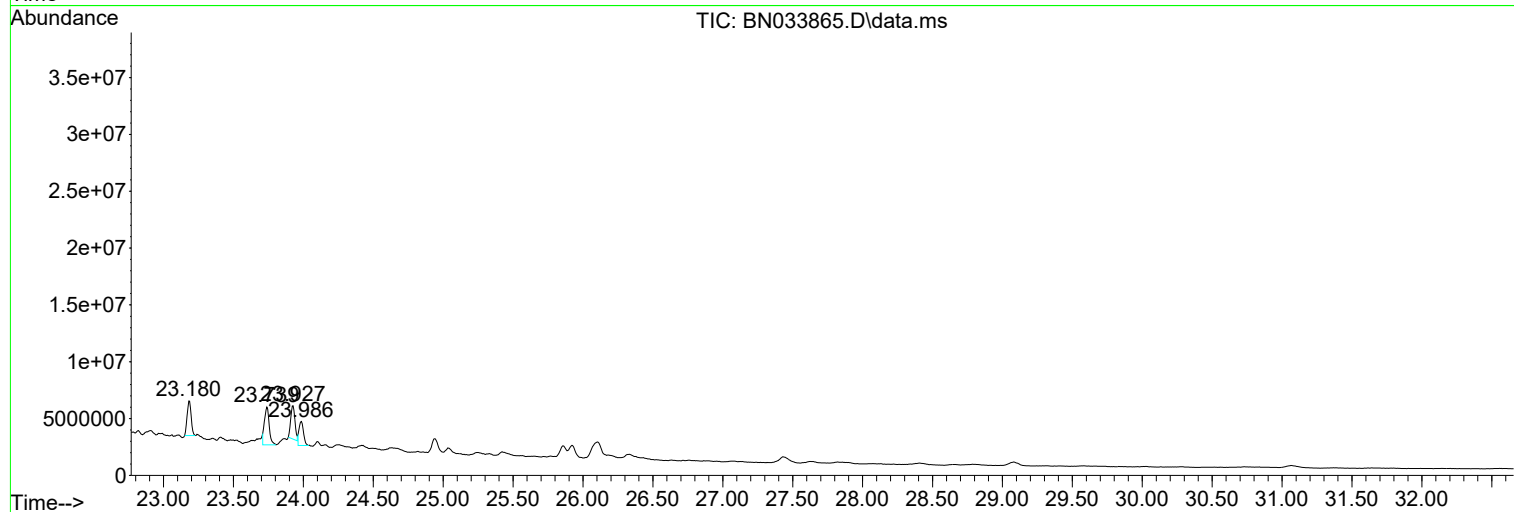
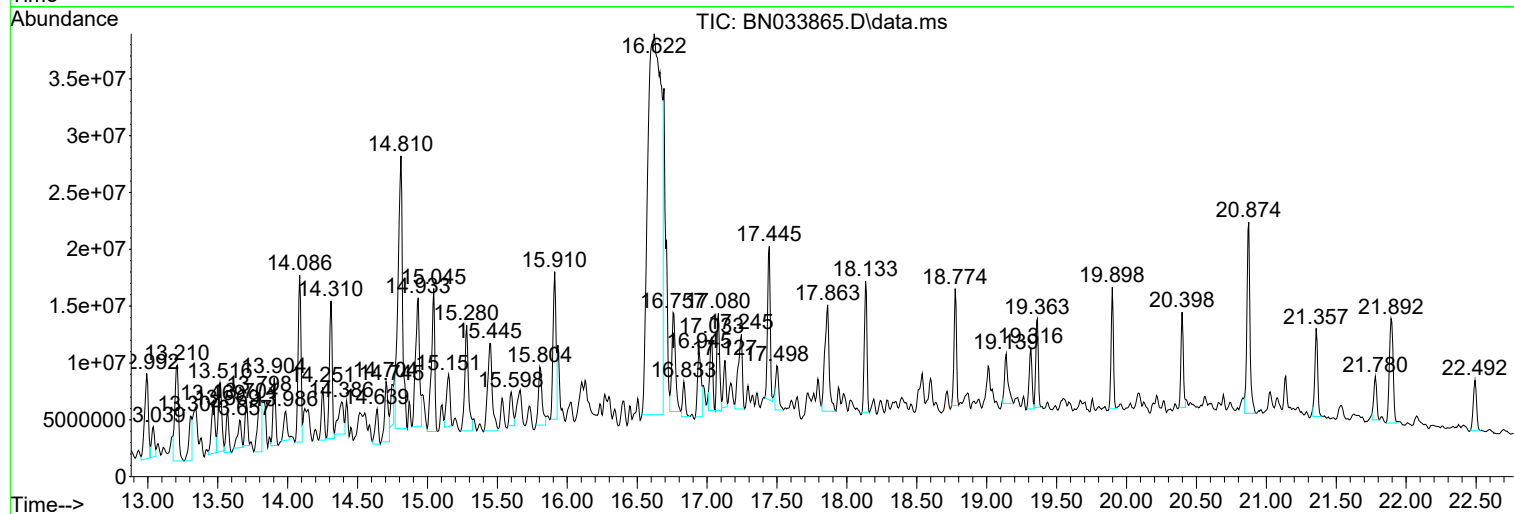
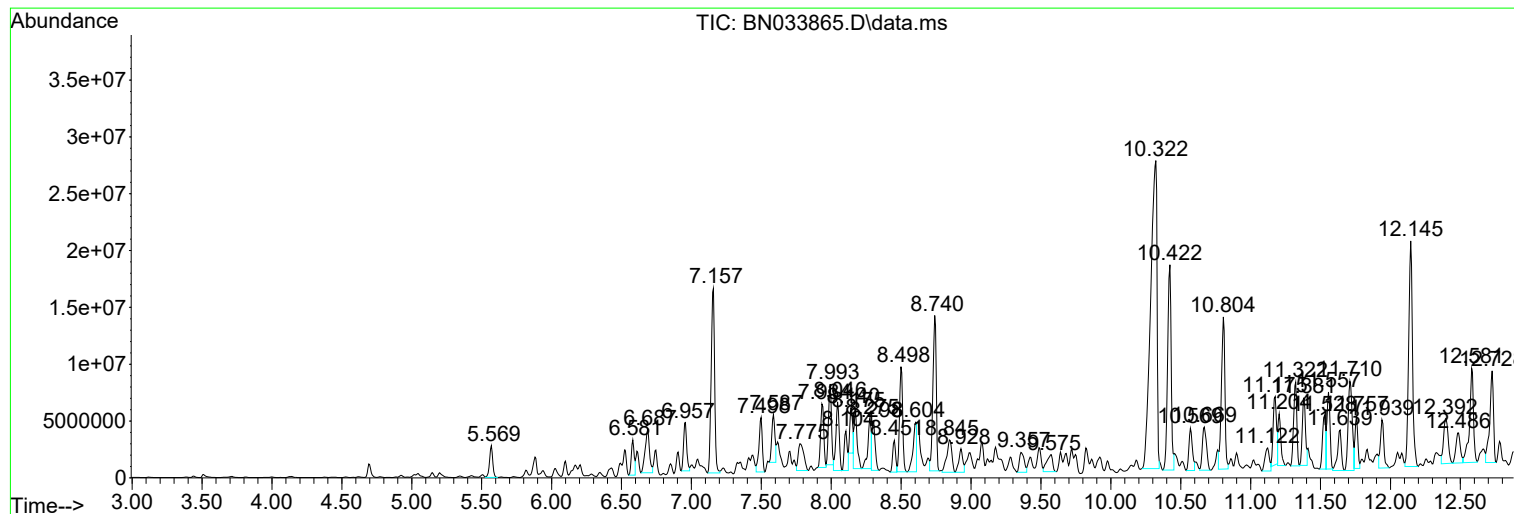
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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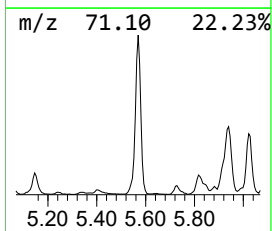
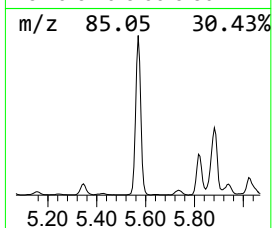
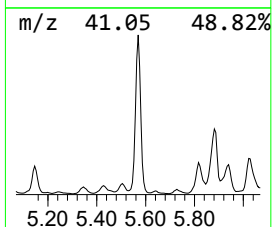
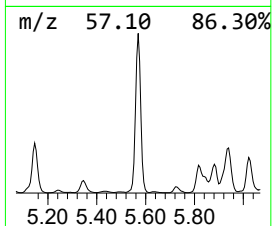
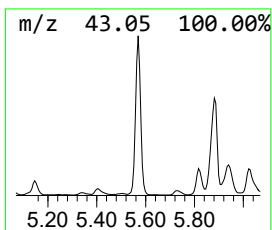
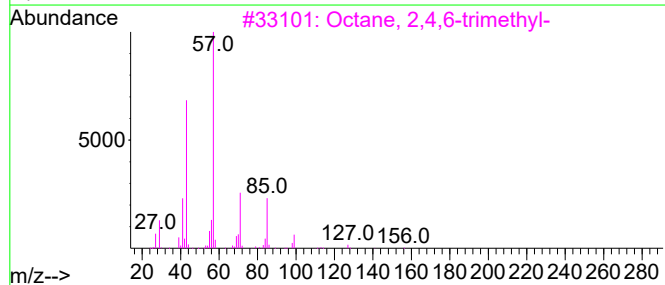
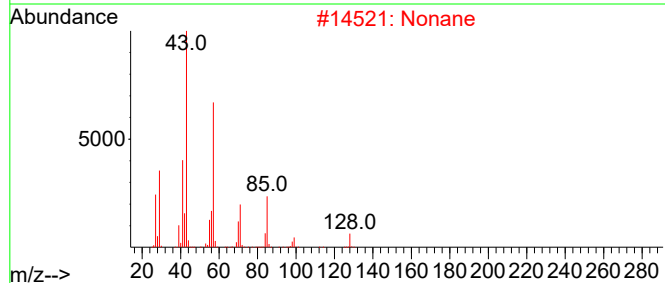
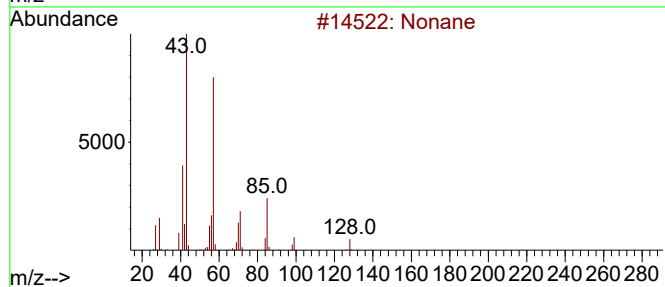
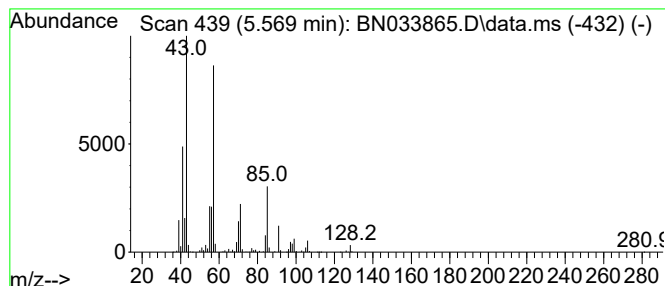
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.569	11.55 ng/ul	4487340	1,4-Dichlorobenzene-d4	7.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane	128	C9H20	000111-84-2	81
2			Nonane	128	C9H20	000111-84-2	72
3			Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	53
4			Nonane	128	C9H20	000111-84-2	52
5			Decane, 2,6,6-trimethyl-	184	C13H28	062108-24-1	50



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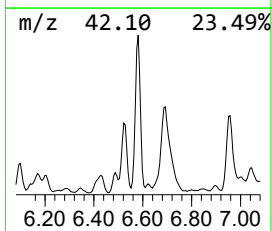
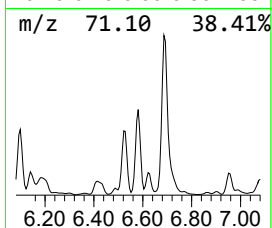
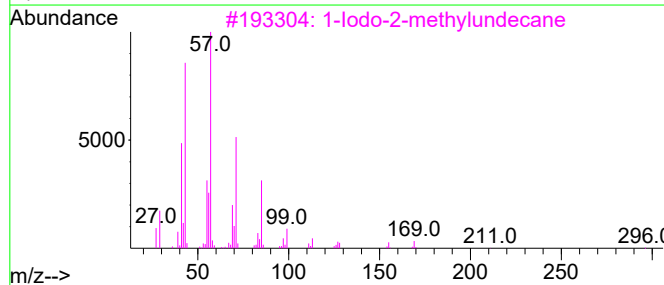
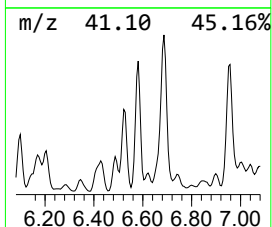
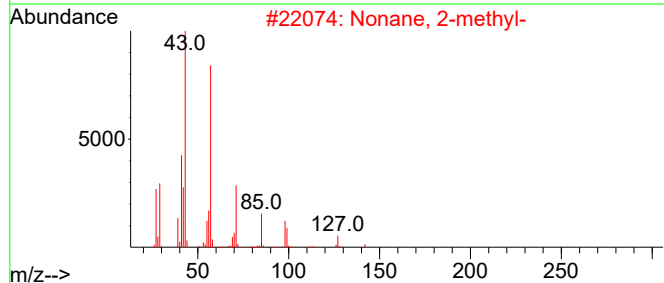
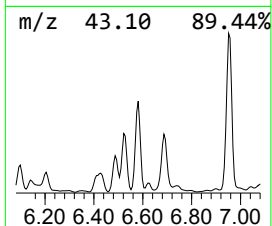
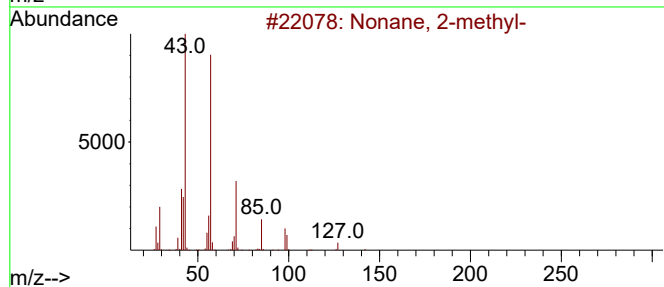
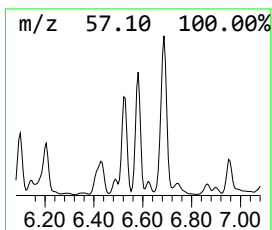
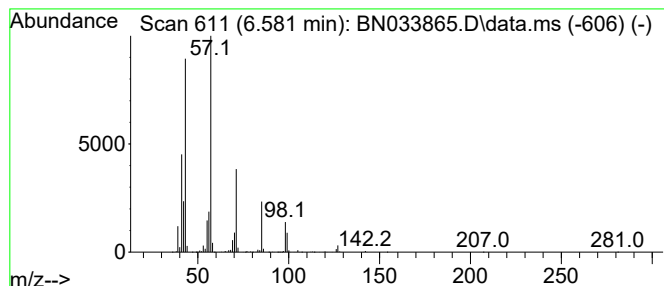
Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD			R.T.
6.581	12.16 ng/ul	4723880	1,4-Dichlorobenzene-d4			7.504
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonane, 2-methyl-		142	C10H22	000871-83-0	93
2	Nonane, 2-methyl-		142	C10H22	000871-83-0	83
3	1-Iodo-2-methylundecane		296	C12H25I	073105-67-6	72
4	Decane, 2-methyl-		156	C11H24	006975-98-0	64
5	Heptane, 2,6-dimethyl-		128	C9H20	001072-05-5	64



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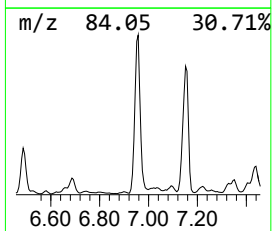
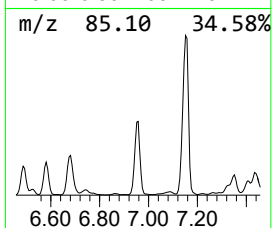
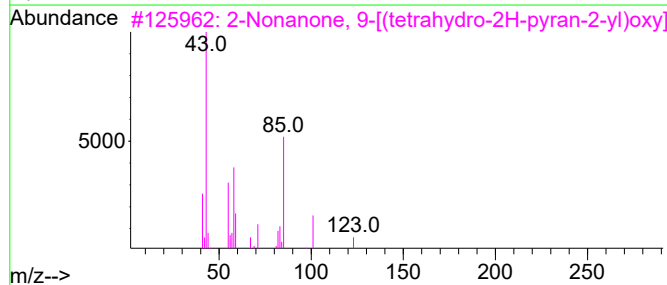
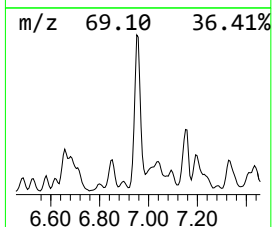
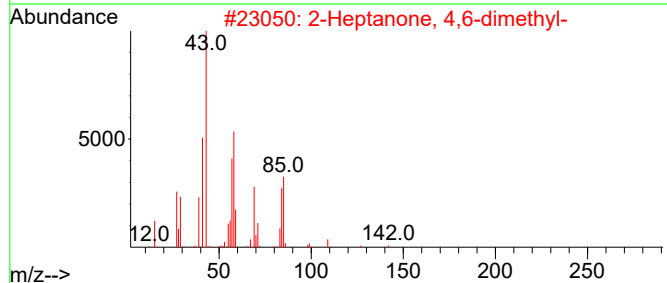
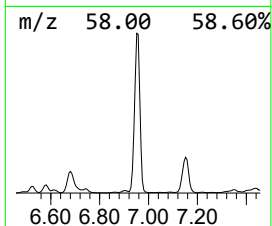
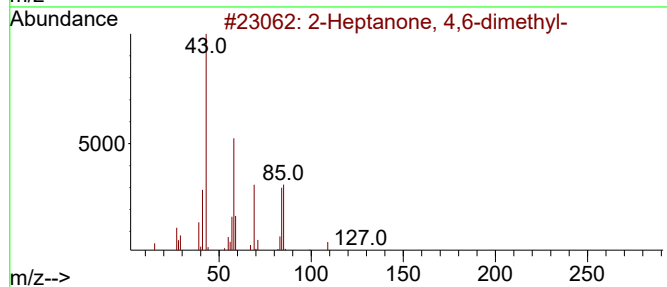
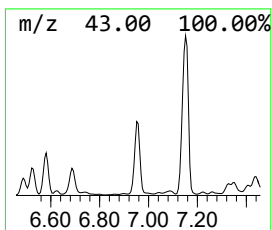
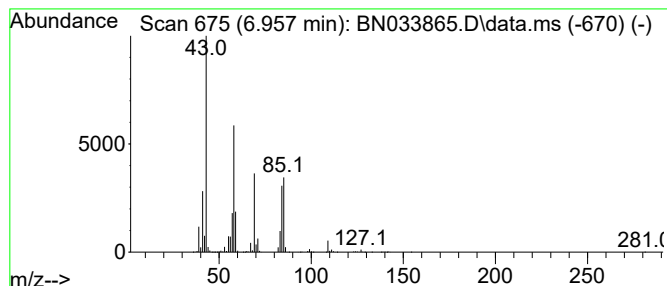
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 2-Heptanone, 4,6-dimethyl- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.957	17.53 ng/ul	6807540	1,4-Dichlorobenzene-d4	7.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	90		
2	2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	78		
3	2-Nonanone, 9-[(tetrahydro-2H-py...	242	C14H26O3	054699-41-1	47		
4	2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	43		
5	2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	43		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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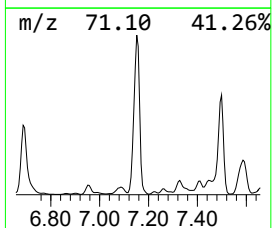
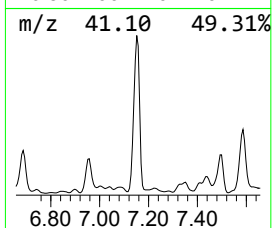
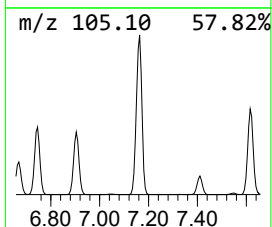
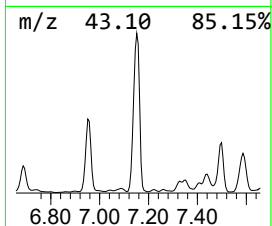
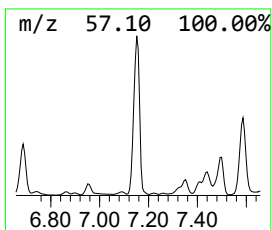
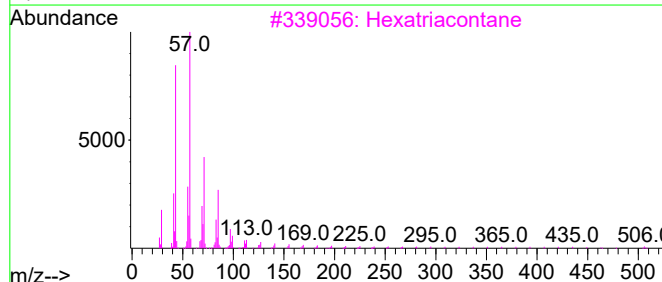
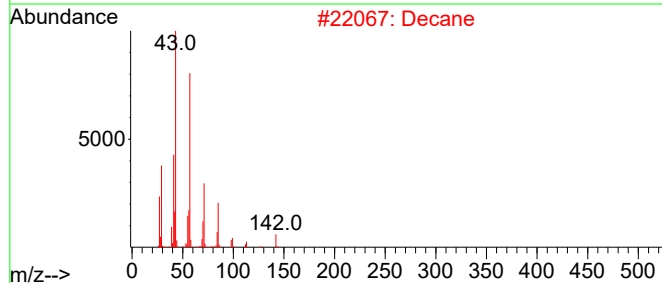
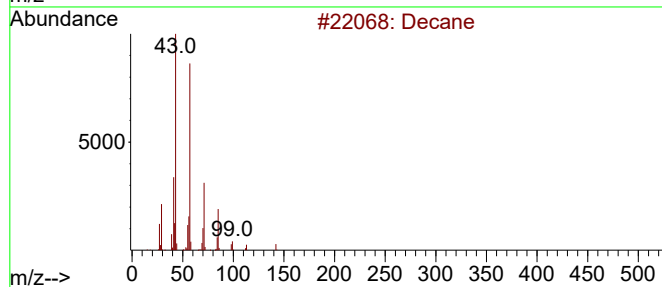
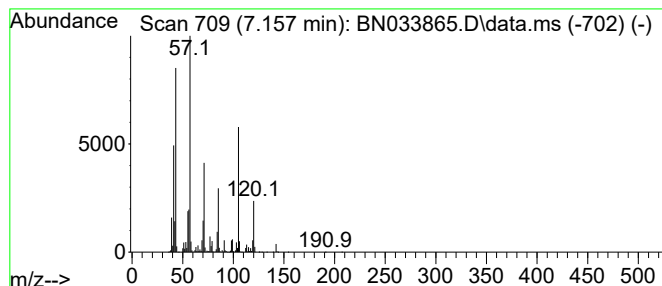
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.157	72.72 ng/ul	28245400	1,4-Dichlorobenzene-d4	7.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane	142	C10H22	000124-18-5	90
2			Decane	142	C10H22	000124-18-5	64
3			Hexatriacontane	507	C36H74	000630-06-8	59
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	56
5			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
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Instrument :
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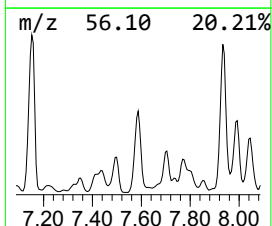
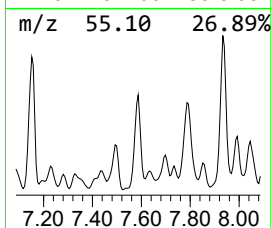
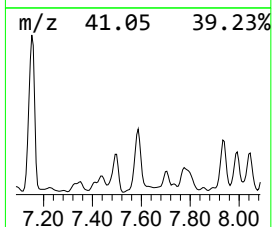
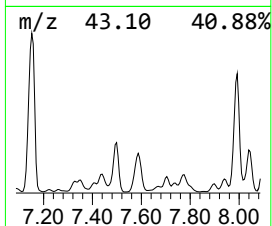
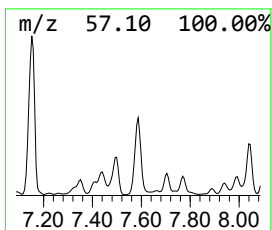
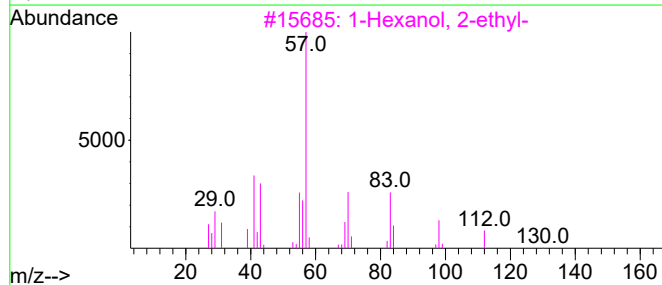
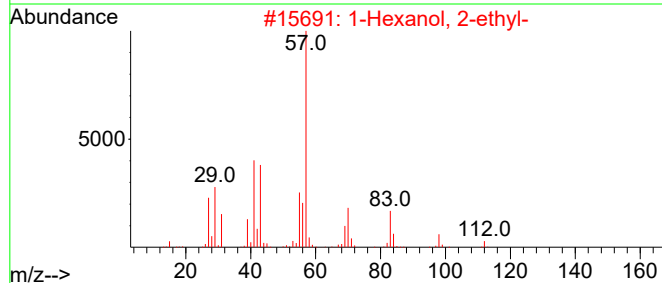
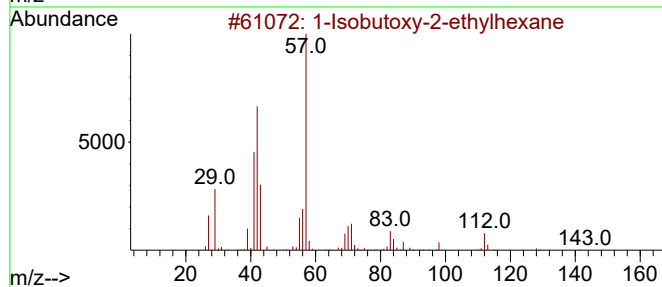
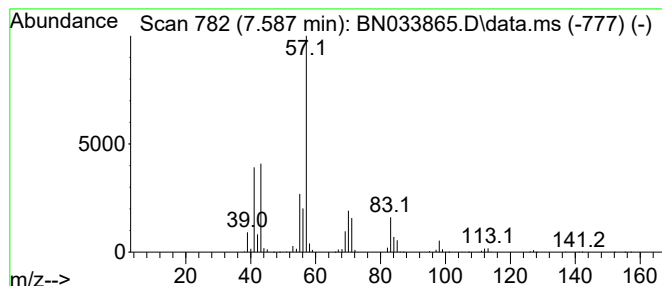
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 1-Isobutoxy-2-ethylhexane Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.587	17.55 ng/ul	6817290	1,4-Dichlorobenzene-d4	7.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Isobutoxy-2-ethylhexane	186	C12H26O	1000139-90-3	74
2			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	72
3			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	72
4			Heptane, 2,3,6-trimethyl-	142	C10H22	004032-93-3	64
5			1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
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Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

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ClientSampleId :
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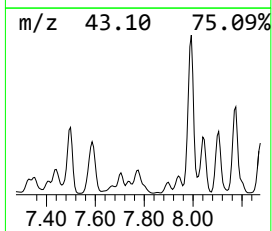
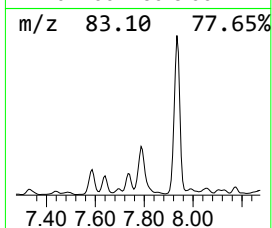
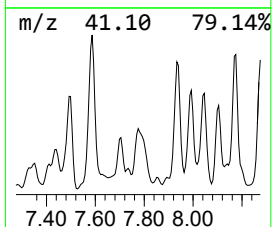
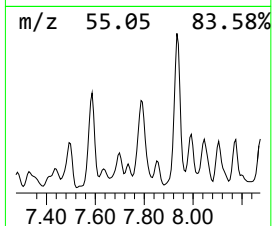
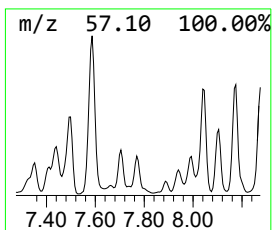
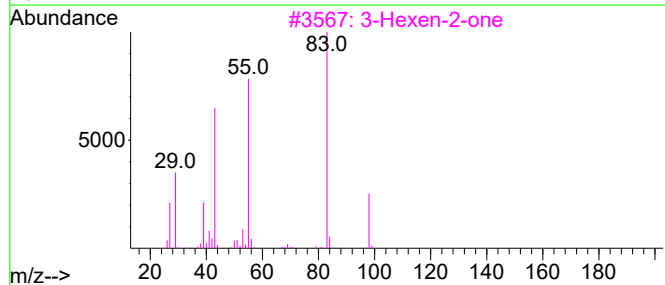
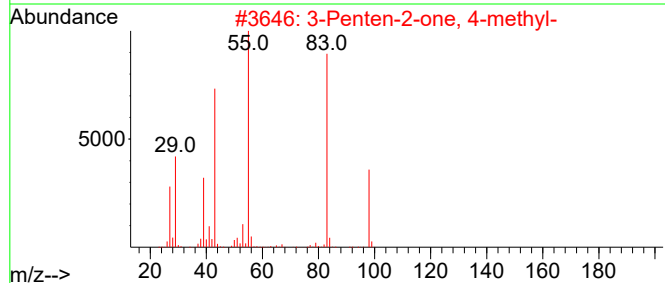
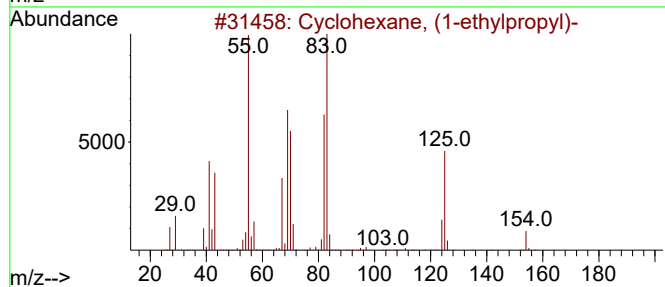
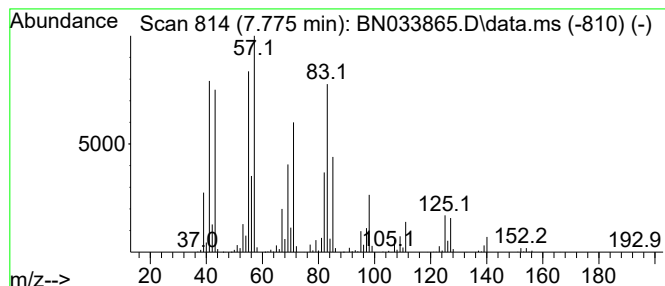
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 (DEL) Alkane: Cyclic7.775 Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.775	16.66 ng/ul	6469150	1,4-Dichlorobenzene-d4	7.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, (1-ethylpropyl)-	154	C11H22	026321-98-2	41
2			3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	30
3			3-Hexen-2-one	98	C6H10O	000763-93-9	30
4			2-Pentene, 3,4-dimethyl-	98	C7H14	024910-63-2	30
5			Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
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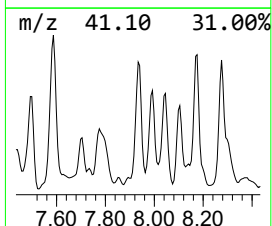
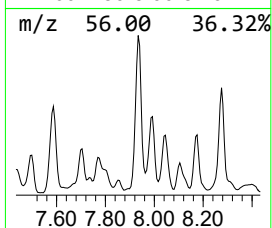
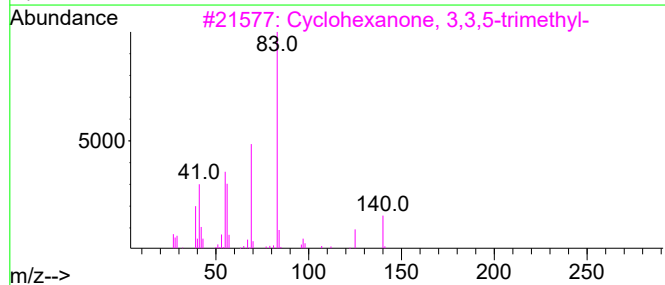
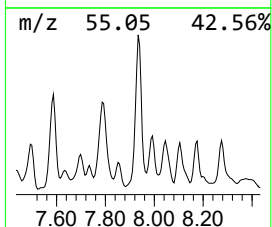
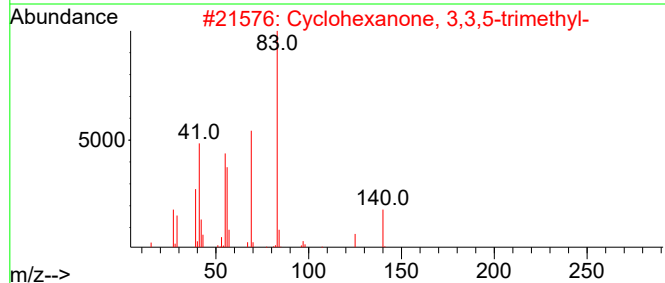
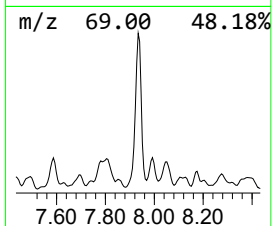
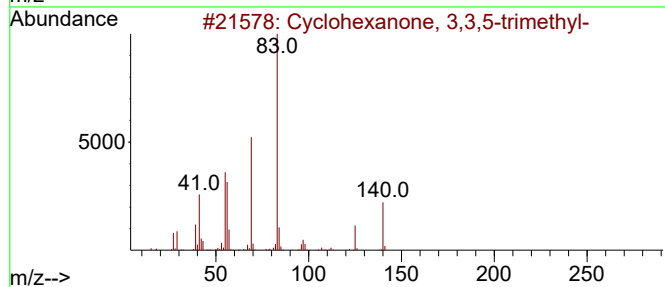
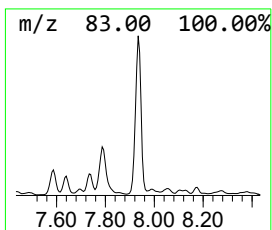
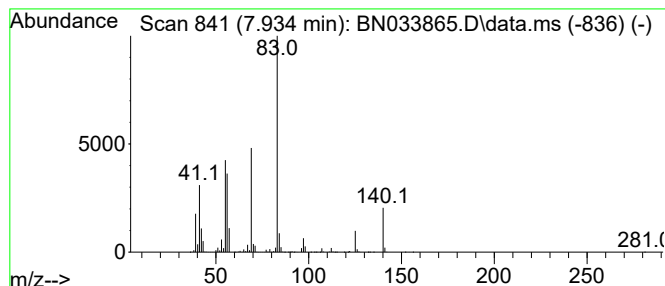
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Cyclohexanone, 3,3,5-trimet... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.934	22.55 ng/ul	8760370	1,4-Dichlorobenzene-d4	7.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	97
2			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	94
3			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	91
4			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	91
5			1,3-Cyclohexanedione, 5,5-dimethyl-	140	C8H12O2	000126-81-8	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
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Acq On : 11 Sep 2024 17:35
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Sample : P3878-18
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ALS Vial : 13 Sample Multiplier: 1

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ClientSampleId :
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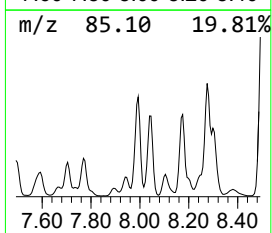
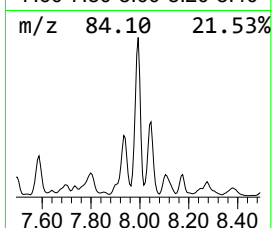
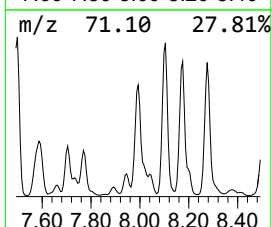
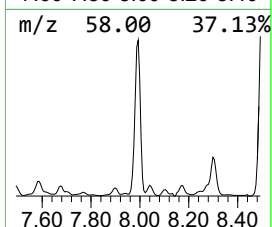
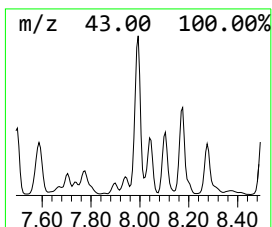
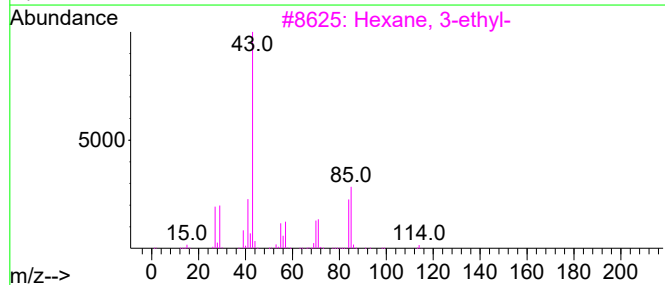
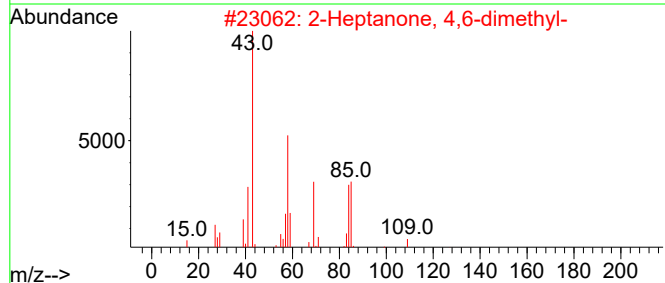
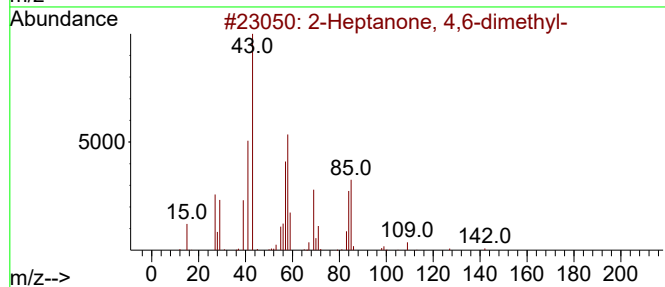
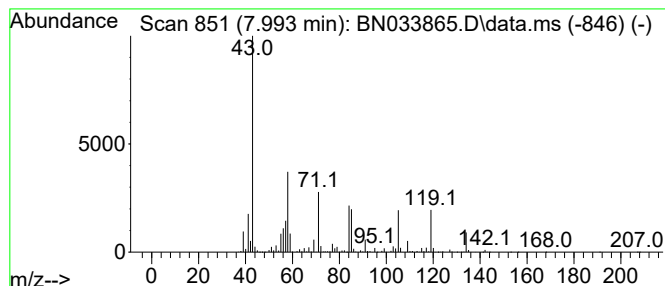
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 unknown-01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.993	26.72 ng/ul	10376700	1,4-Dichlorobenzene-d4	7.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	35
2			2-Heptanone, 4,6-dimethyl-	142	C9H18O	019549-80-5	32
3			Hexane, 3-ethyl-	114	C8H18	000619-99-8	27
4			Hexane, 3-ethyl-	114	C8H18	000619-99-8	22
5			Heptane, 2,3-dimethyl-	128	C9H20	003074-71-3	16



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
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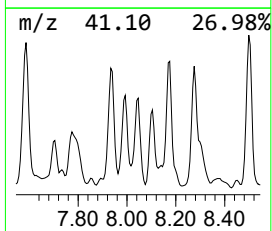
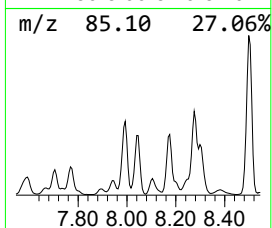
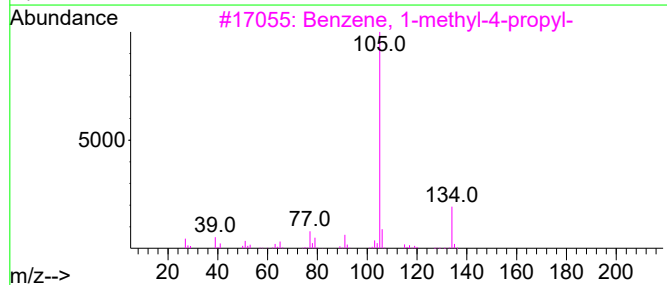
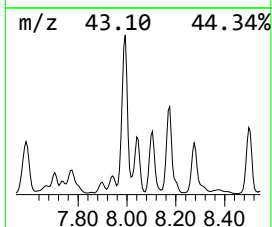
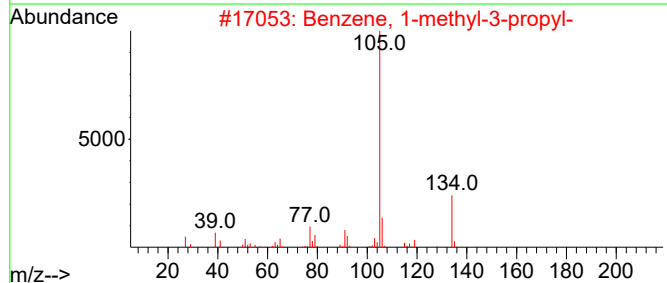
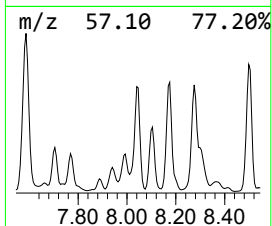
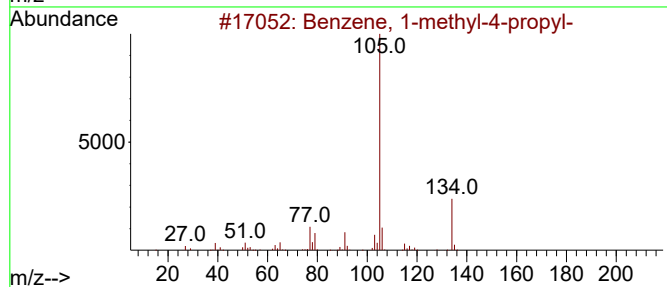
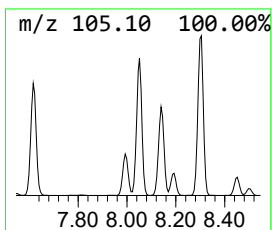
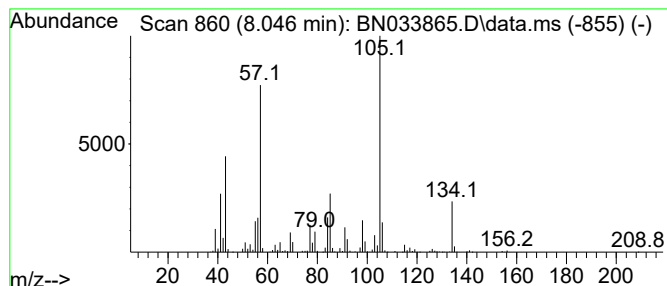
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzene, 1-methyl-4-propyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.046	25.24 ng/ul	9804530	1,4-Dichlorobenzene-d4	7.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	83
2			Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	60
3			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	46
4			Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	46
5			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
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Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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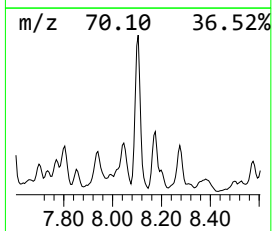
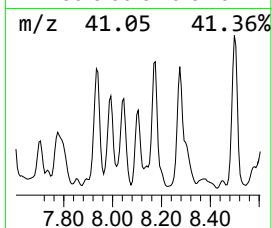
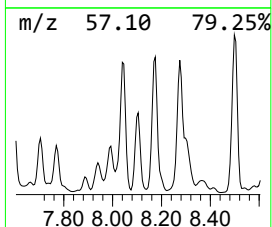
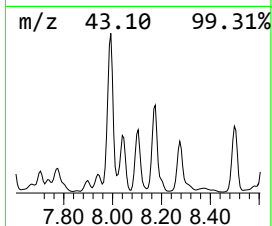
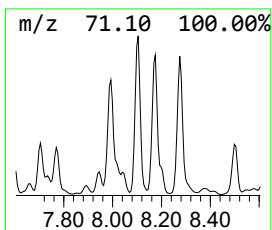
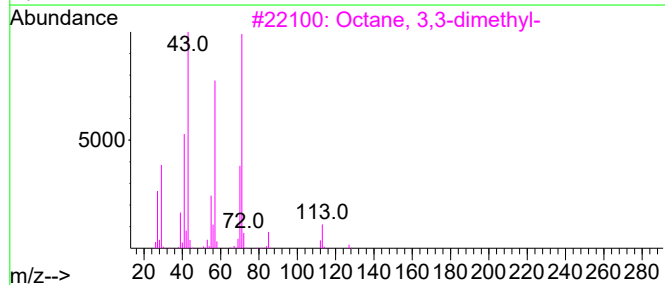
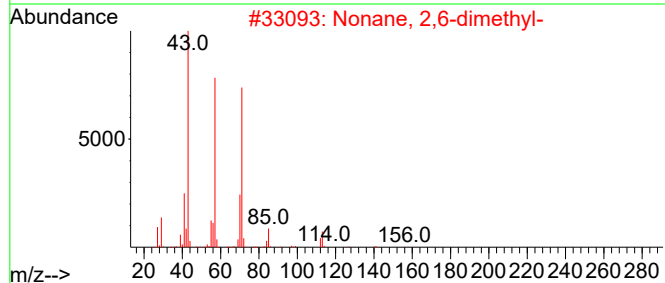
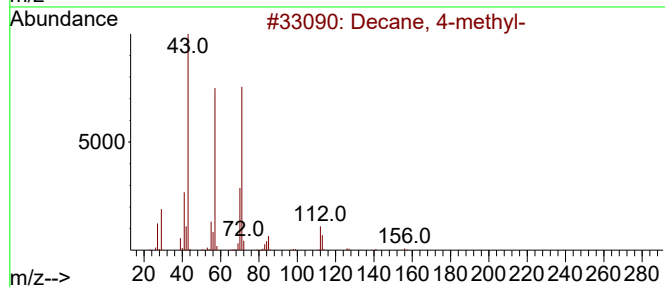
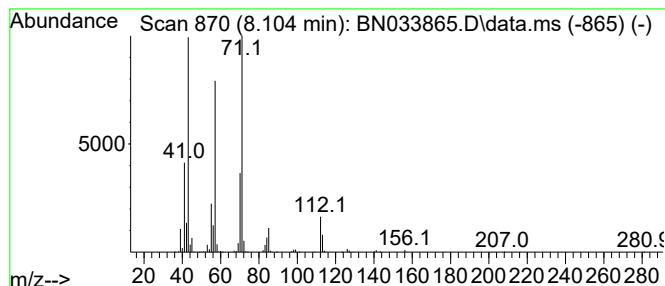
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.104	13.78 ng/ul	5353230	1,4-Dichlorobenzene-d4	7.504

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Decane, 4-methyl-	156	C11H24	002847-72-5	97
2		Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	91
3		Octane, 3,3-dimethyl-	142	C10H22	004110-44-5	86
4		Decane, 4-methyl-	156	C11H24	002847-72-5	76
5		Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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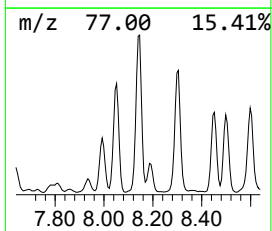
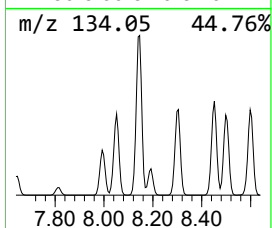
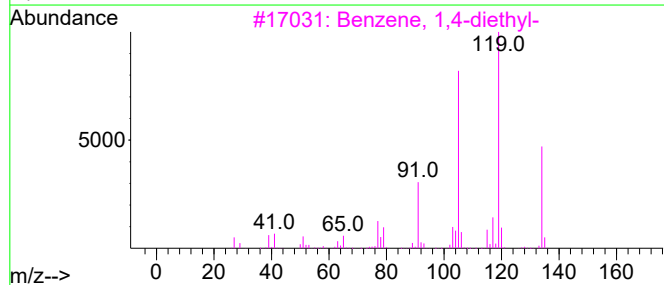
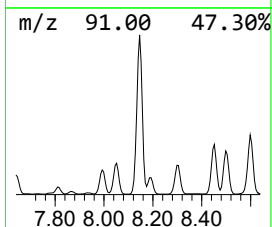
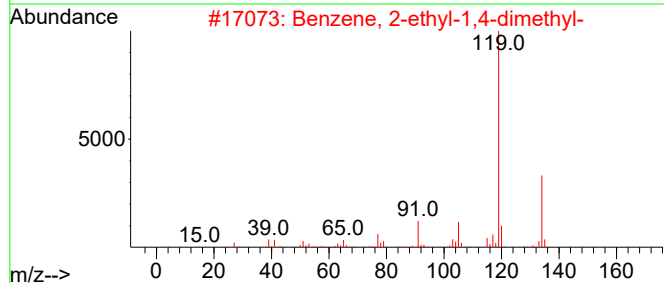
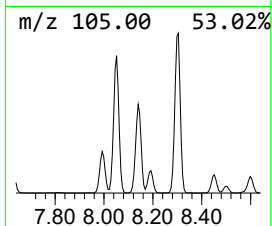
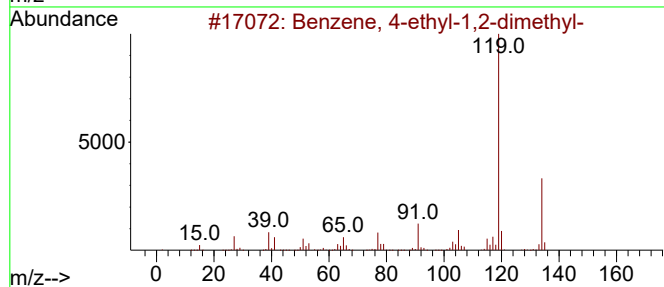
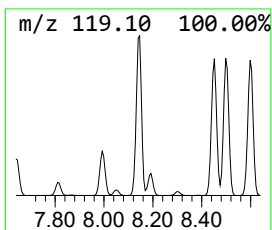
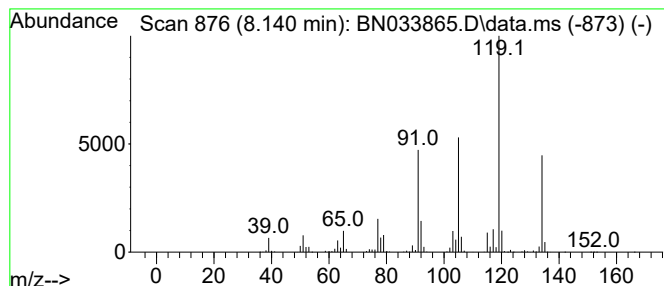
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.140	14.59 ng/ul	5665160	1,4-Dichlorobenzene-d4	7.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
3			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	92
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
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Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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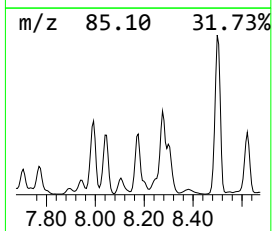
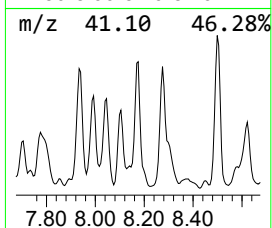
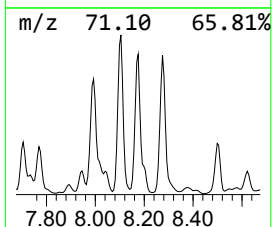
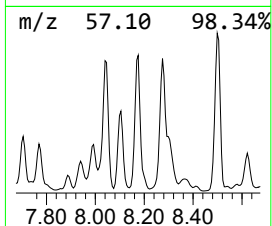
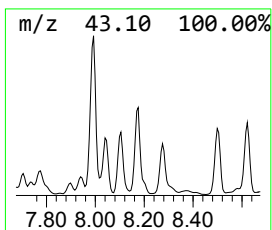
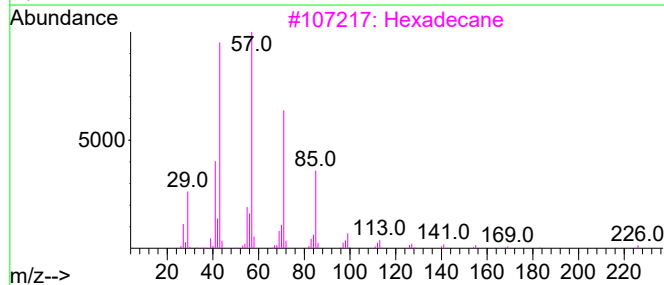
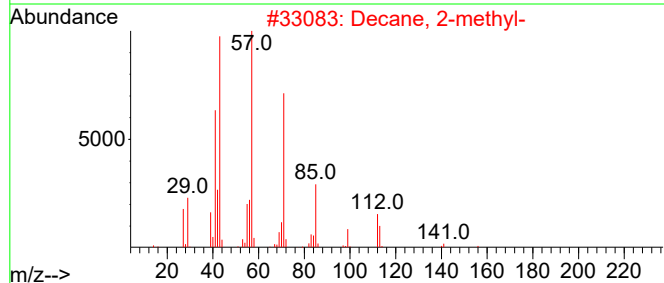
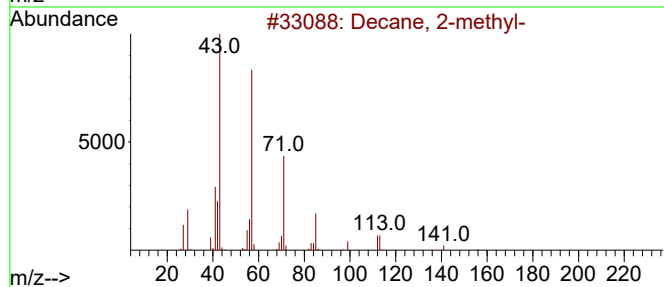
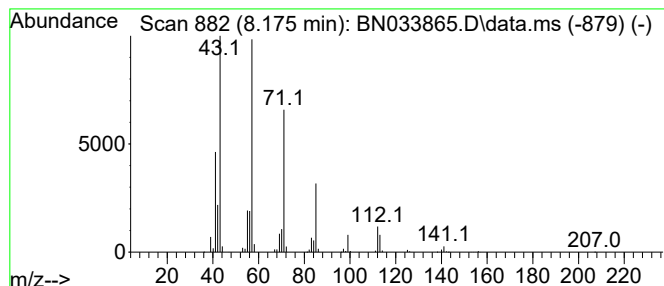
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.175	20.19 ng/ul	7843930	1,4-Dichlorobenzene-d4	7.504

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Decane, 2-methyl-	156	C11H24	006975-98-0	90
2		Decane, 2-methyl-	156	C11H24	006975-98-0	90
3		Hexadecane	226	C16H34	000544-76-3	72
4		Tetradecane	198	C14H30	000629-59-4	72
5		Heptadecane	240	C17H36	000629-78-7	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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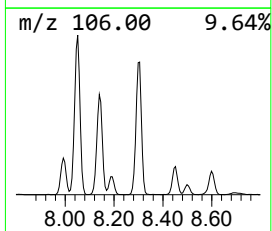
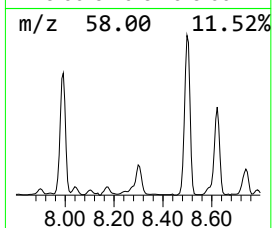
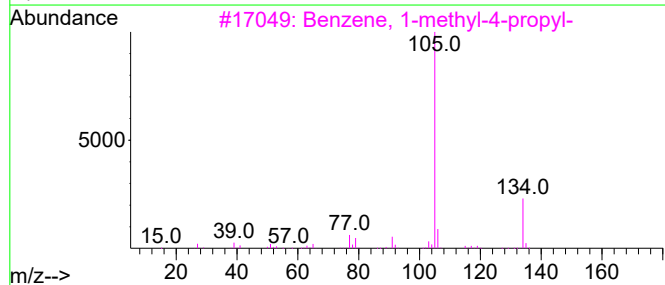
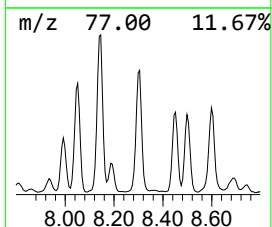
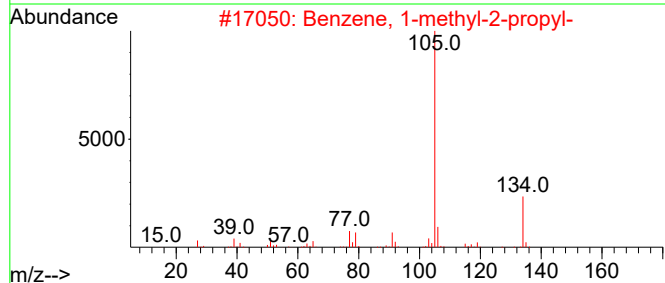
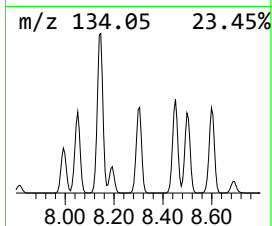
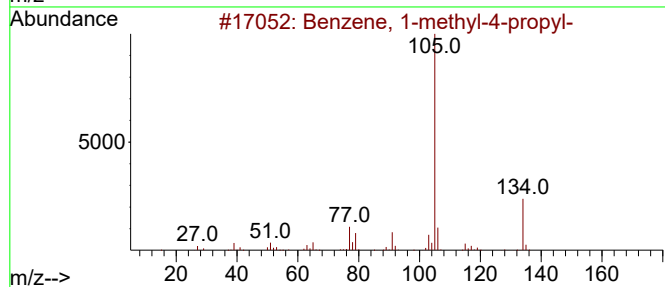
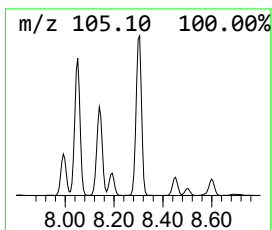
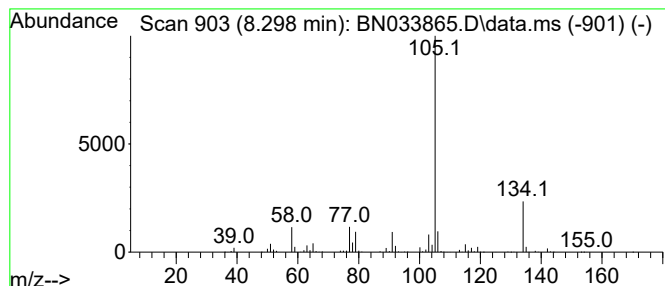
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Benzene, 1-methyl-2-propyl- Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.298	16.98 ng/ul	6595080	1,4-Dichlorobenzene-d4	7.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	94
2			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	93
3			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	87
4			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	87
5			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
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Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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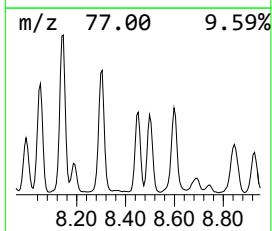
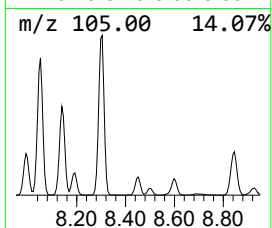
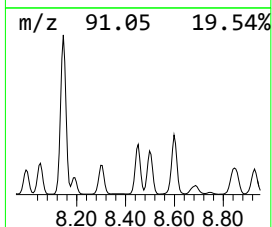
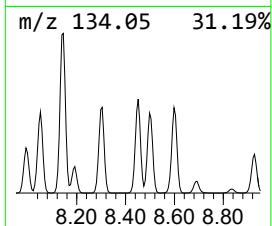
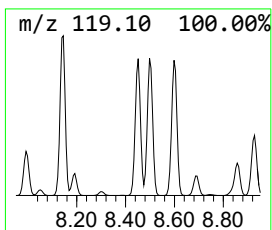
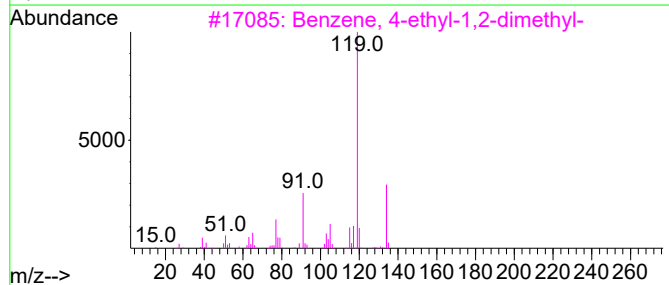
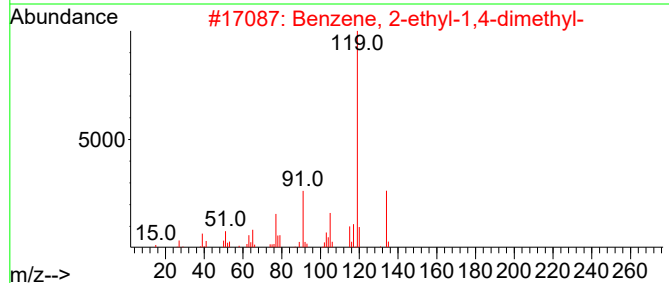
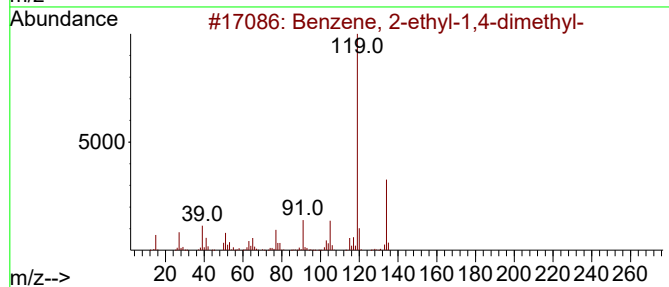
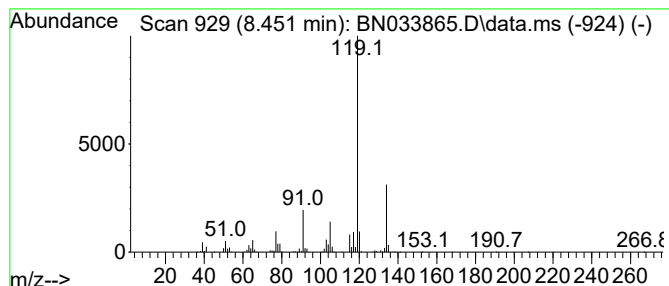
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.451	10.31 ng/ul	4004300	1,4-Dichlorobenzene-d4	7.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
5			o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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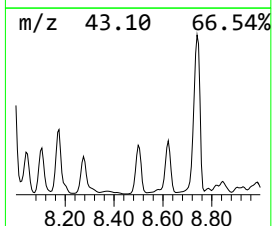
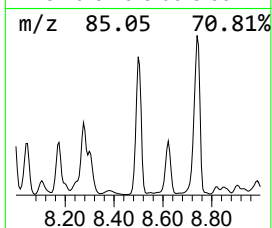
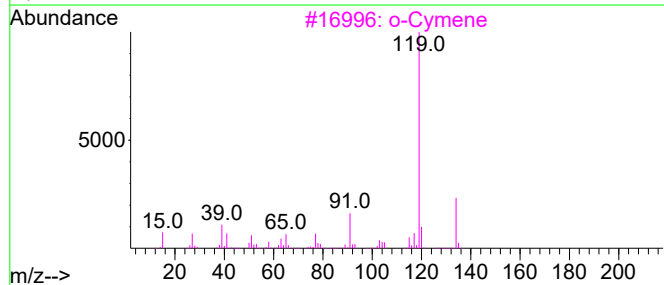
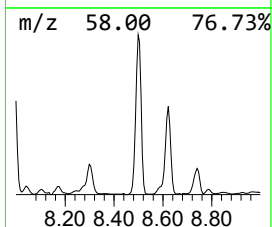
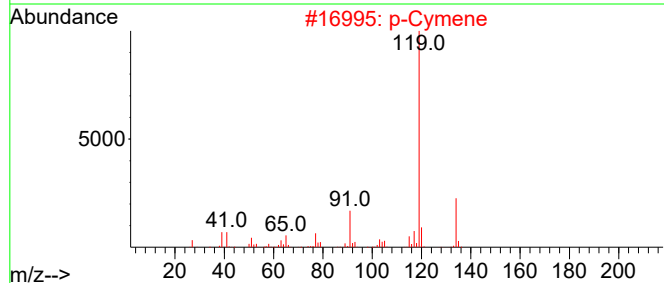
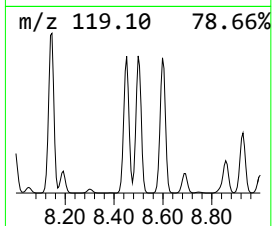
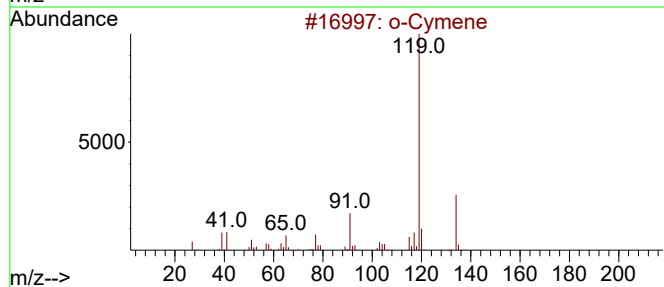
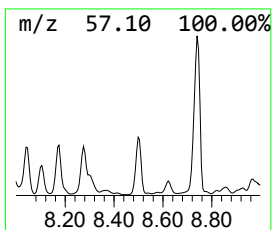
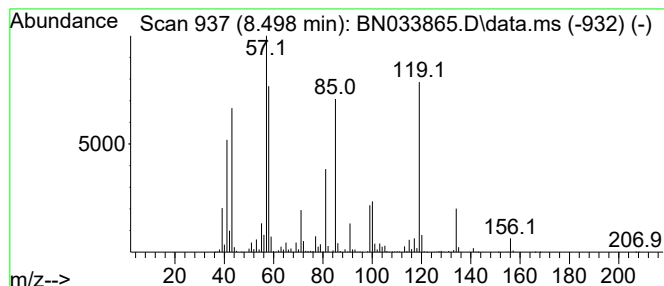
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 o-Cymene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.498	36.47 ng/ul	14165500	1,4-Dichlorobenzene-d4	7.504

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		o-Cymene	134	C10H14	000527-84-4	68
2		p-Cymene	134	C10H14	000099-87-6	68
3		o-Cymene	134	C10H14	000527-84-4	68
4		o-Cymene	134	C10H14	000527-84-4	68
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
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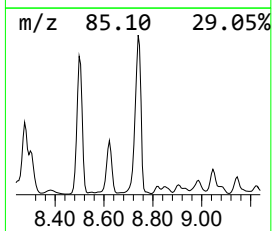
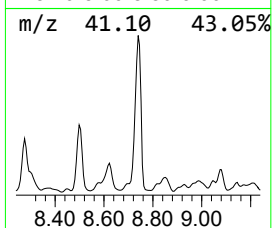
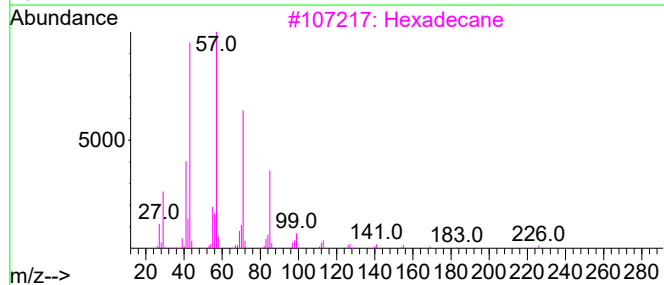
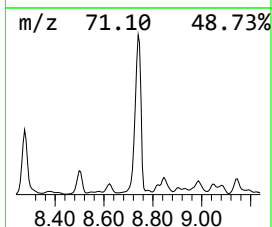
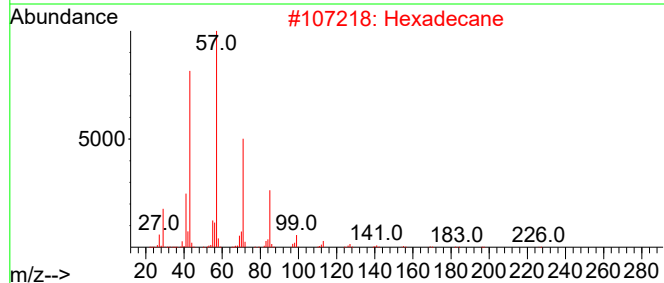
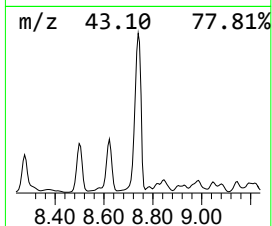
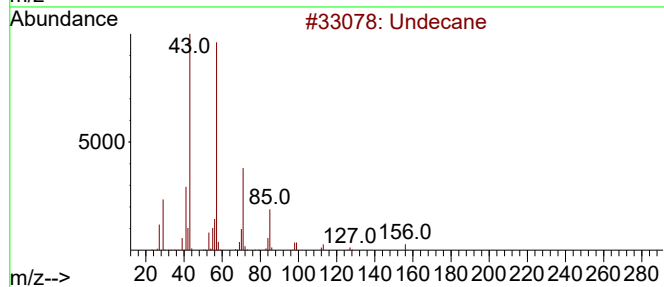
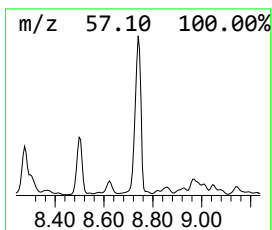
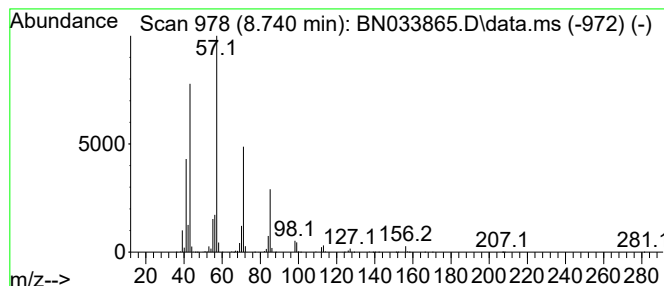
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.740	57.94 ng/ul	22506500	1,4-Dichlorobenzene-d4	7.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane	156	C11H24	001120-21-4	94
2			Hexadecane	226	C16H34	000544-76-3	90
3			Hexadecane	226	C16H34	000544-76-3	90
4			Decane	142	C10H22	000124-18-5	90
5			Carbonic acid, prop-1-en-2-yl te...	298	C18H34O3	1000382-90-9	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
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Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

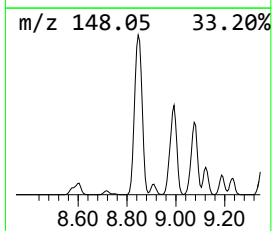
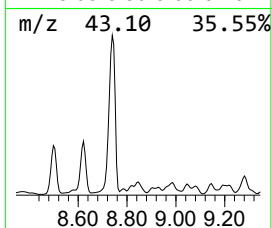
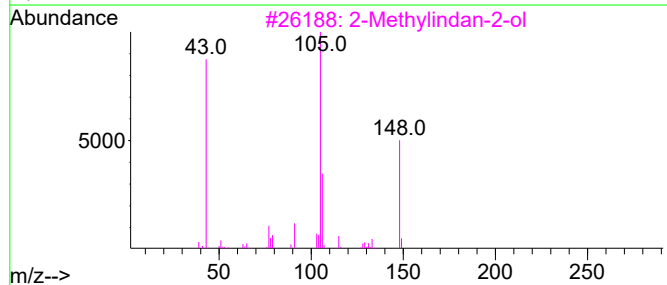
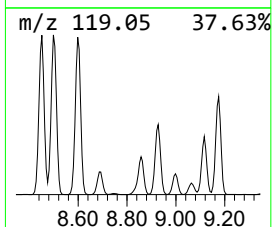
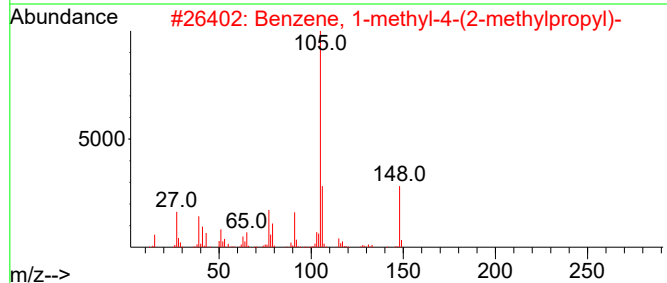
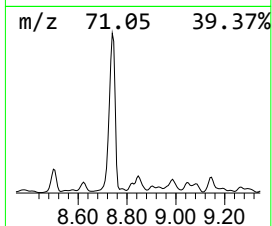
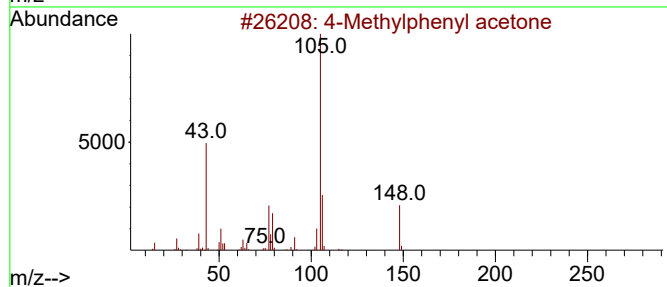
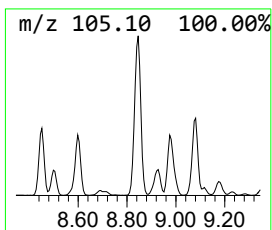
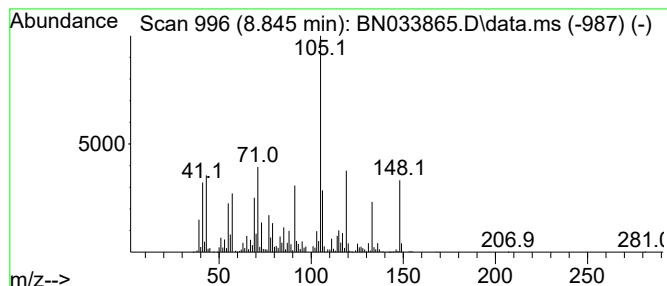
Instrument :
BNA_N
ClientSampleId :
DDCC1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 unknown-02 Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD			R.T.
8.845	18.83 ng/ul	7312680	1,4-Dichlorobenzene-d4			7.504
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	4-Methylphenyl acetone		148	C10H12O	002096-86-8	42
2	Benzene, 1-methyl-4-(2-methylpro...		148	C11H16	005161-04-6	38
3	2-Methylindan-2-ol		148	C10H12O	033223-84-6	27
4	Benzeneethanol, .beta.-methyl-		136	C9H12O	001123-85-9	25
5	3-Buten-2-ol, 4-phenyl-		148	C10H12O	017488-65-2	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

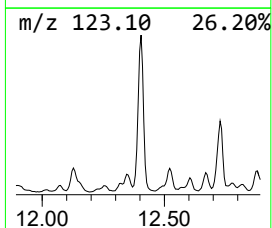
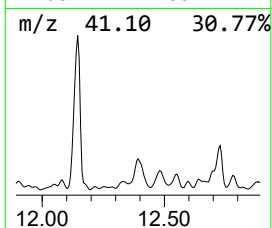
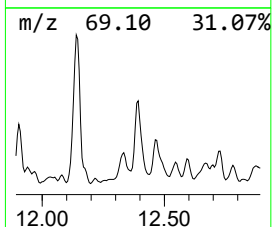
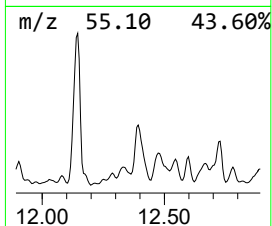
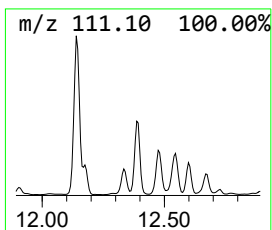
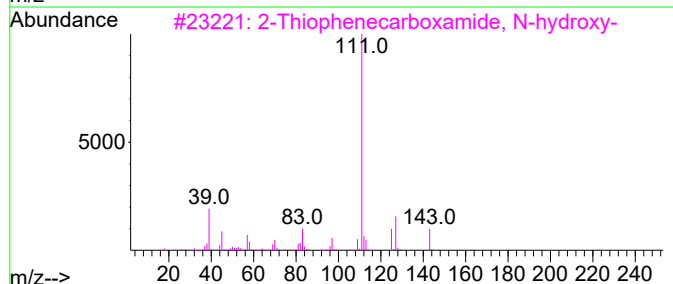
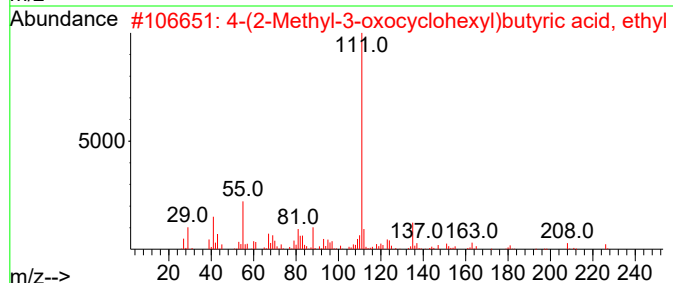
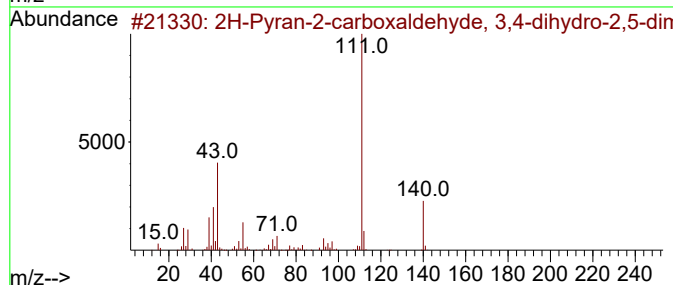
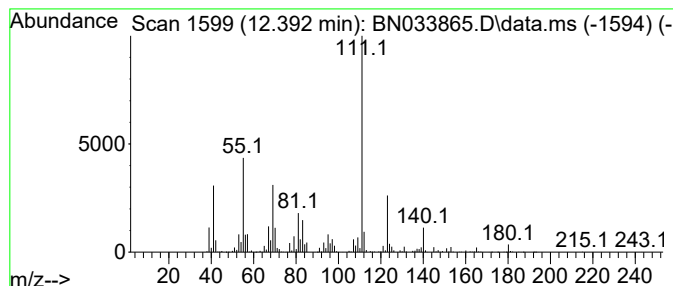
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 24 2H-Pyran-2-carboxaldehyde, ... Concentration Rank 48

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.392	8.53 ng/ul	8983920	Acenaphthene-d10	14.145	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Pyran-2-carboxaldehyde, 3,4-d...	140	C8H12O2	001920-21-4	53
2	4-(2-Methyl-3-oxocyclohexyl)buty...	226	C13H22O3	092486-17-4	53
3	2-Thiophenecarboxamide, N-hydroxy-	143	C5H5NO2S	017698-15-6	50
4	Cyclohexane, 1,4-dimethyl-2-octa...	364	C26H52	055282-02-5	50
5	4-Butyl-5-methylpyrazalone	154	C8H14N2O	1010289-28-0	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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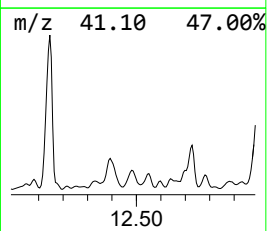
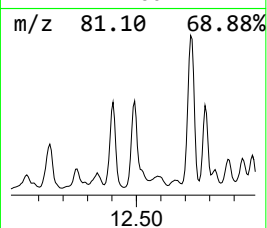
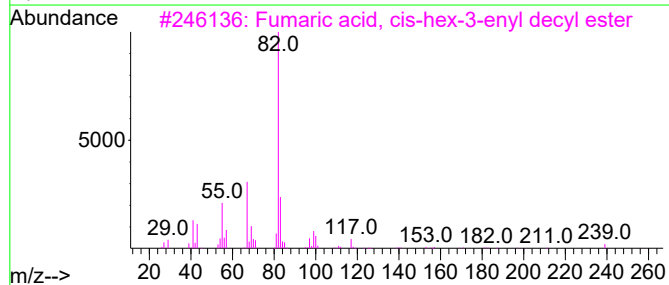
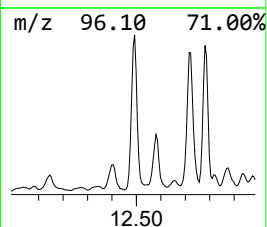
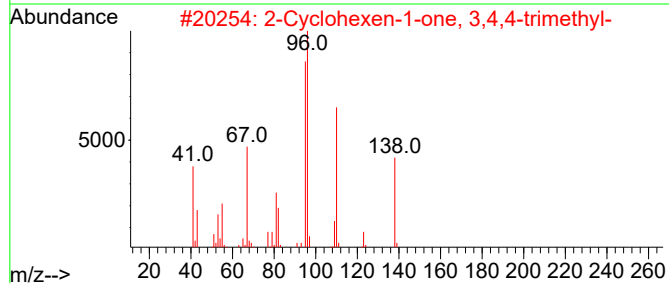
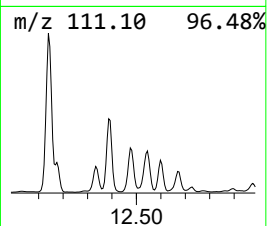
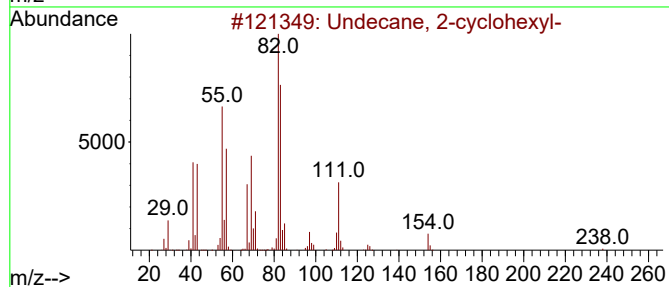
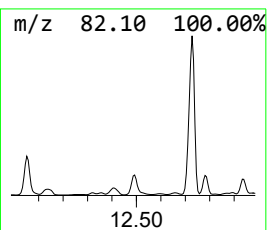
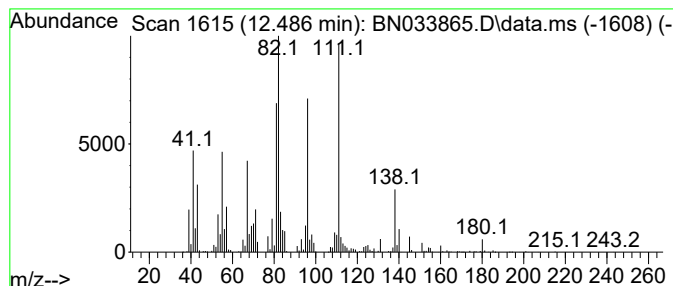
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 25 (DEL) Alkane: Cyclic12.486 Concentration Rank 55

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.486	5.49 ng/ul	5775650	Acenaphthene-d10	14.145

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 2-cyclohexyl-	238	C17H34	013151-77-4	35
2			2-Cyclohexen-1-one, 3,4,4-trimet...	138	C9H14O	017299-41-1	30
3			Fumaric acid, cis-hex-3-enyl dec...	338	C20H34O4	1010348-86-6	27
4			Octane, 2-cyclohexyl-	196	C14H28	002883-05-8	27
5			2(1H)-Pyridinethione	111	C5H5NS	002637-34-5	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

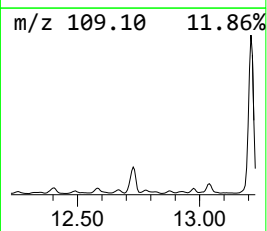
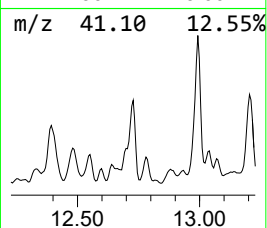
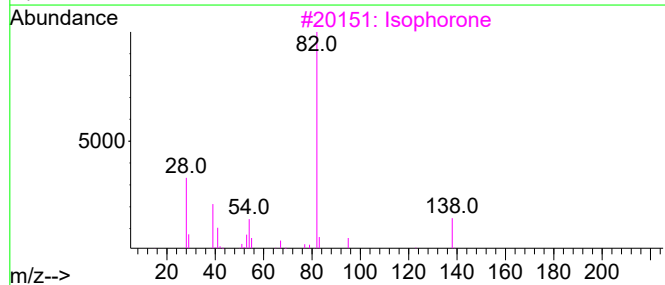
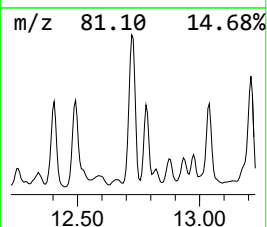
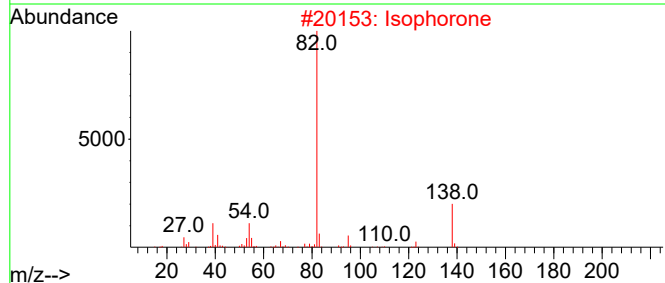
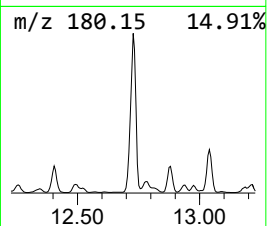
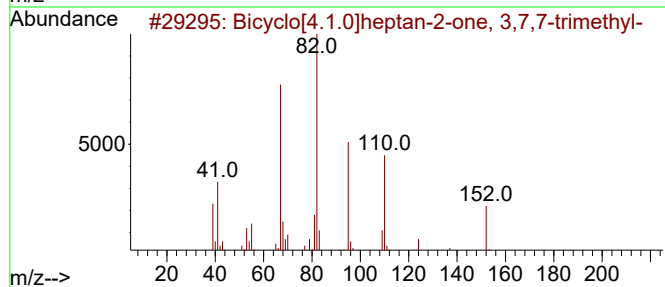
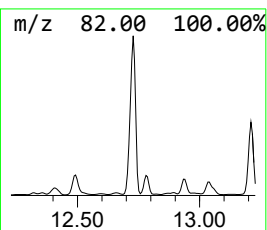
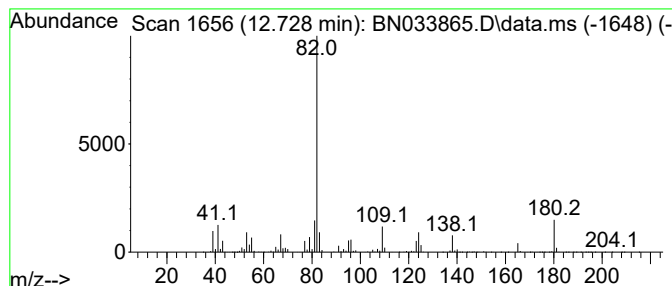
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 26 Bicyclo[4.1.0]heptan-2-one,... Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.728	14.04 ng/ul	14783500	Acenaphthene-d10		14.145	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Bicyclo[4.1.0]heptan-2-one, 3,7,...		152	C10H16O	000497-62-1	64
2	Isophorone		138	C9H14O	000078-59-1	53
3	Isophorone		138	C9H14O	000078-59-1	52
4	Isophorone		138	C9H14O	000078-59-1	52
5	cis-Hexenyl octyne carbonate		236	C15H24O2	068480-29-5	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
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Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

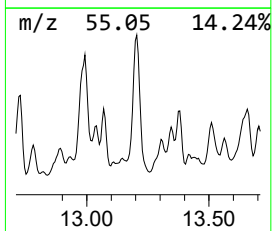
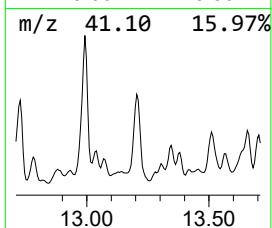
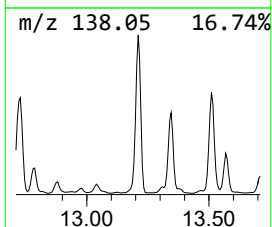
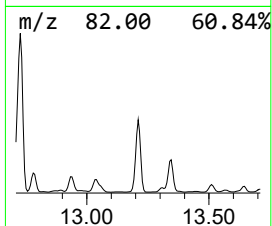
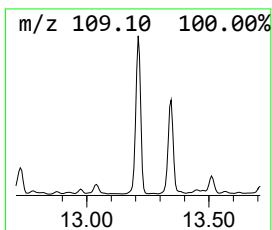
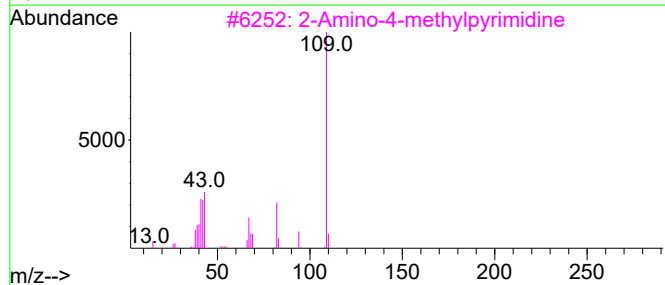
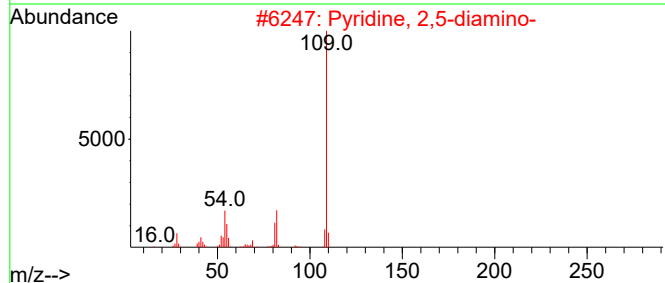
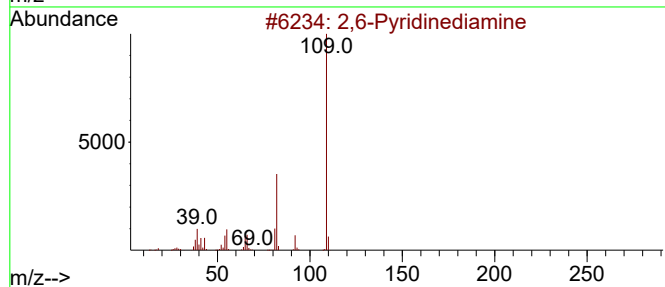
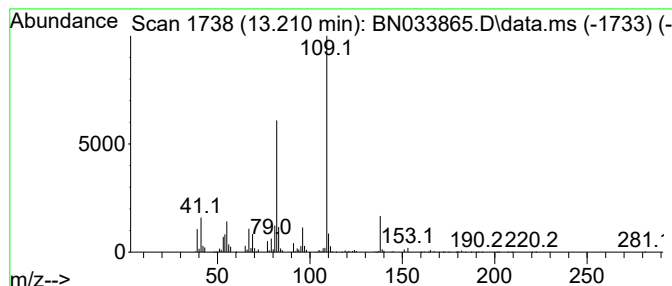
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 28 2,6-Pyridinediamine Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD			R.T.
13.210	14.55 ng/ul	15313400	Acenaphthene-d10			14.145
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2,6-Pyridinediamine		109	C5H7N3	000141-86-6	64
2	Pyridine, 2,5-diamino-		109	C5H7N3	004318-76-7	58
3	2-Amino-4-methylpyrimidine		109	C5H7N3	000108-52-1	47
4	2,6-Pyridinediamine		109	C5H7N3	000141-86-6	47
5	Phenol, 4-amino-		109	C6H7NO	000123-30-8	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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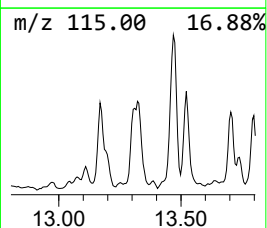
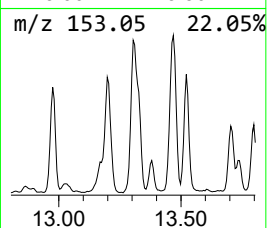
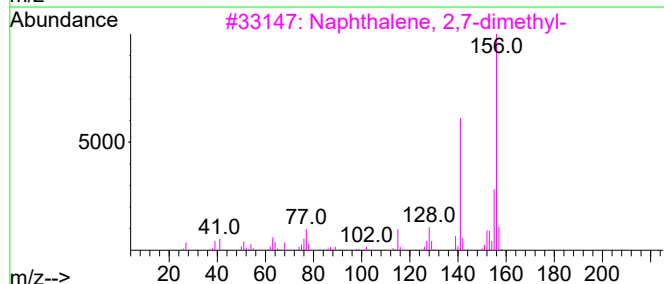
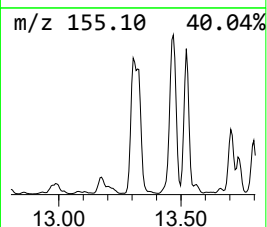
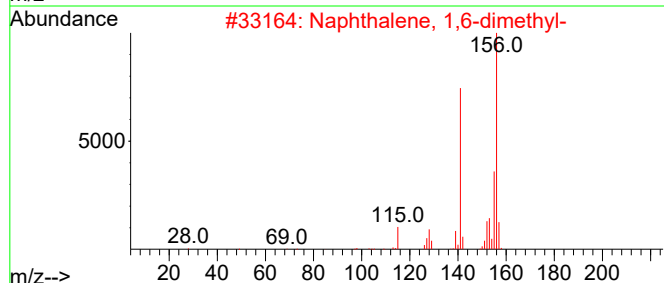
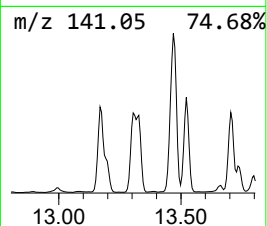
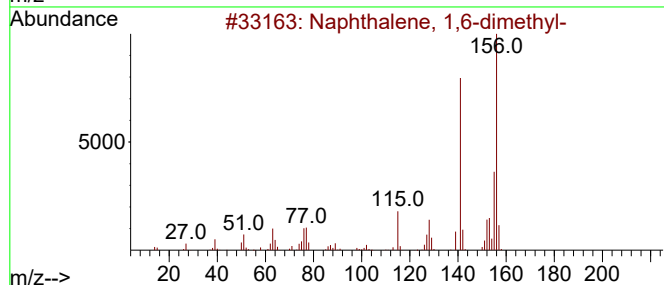
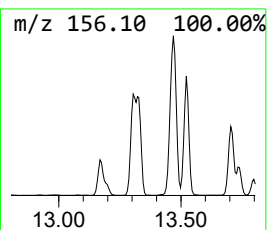
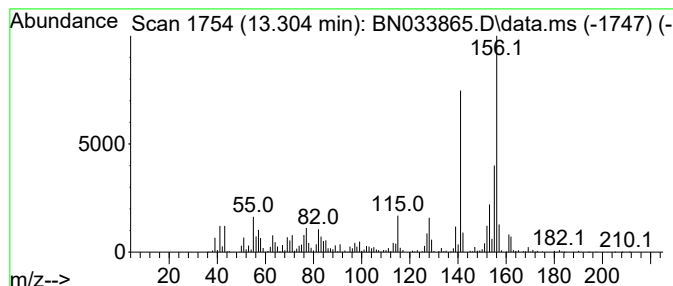
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 29 Naphthalene, 1,6-dimethyl- Concentration Rank 53

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.304	6.21 ng/ul	6534760	Acenaphthene-d10	14.145

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
2			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
3			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
5			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
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Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

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ClientSampleId :
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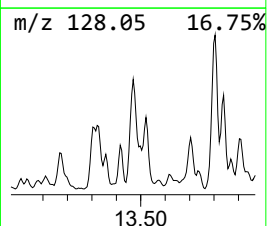
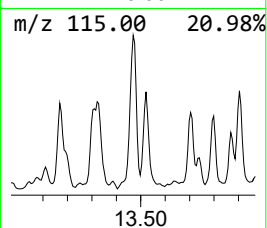
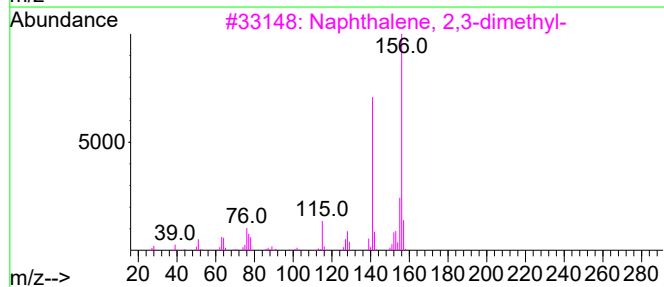
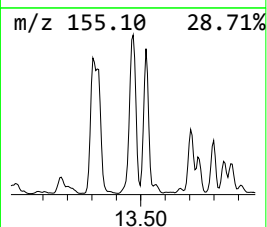
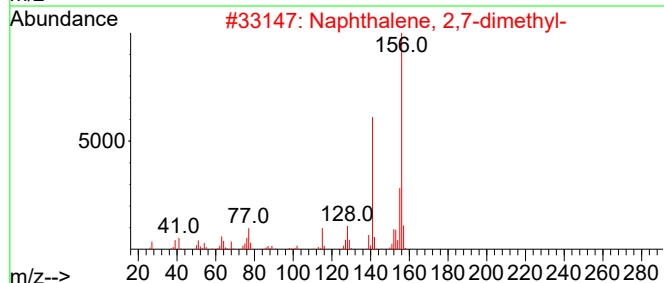
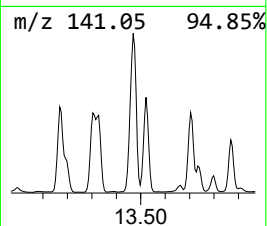
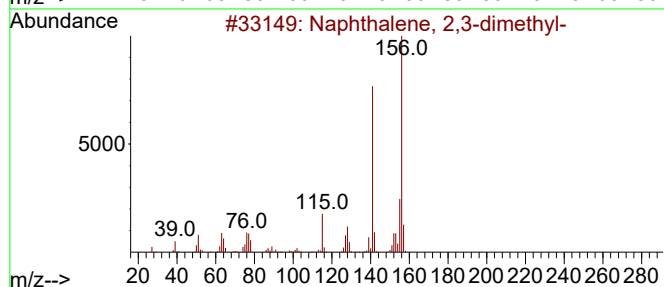
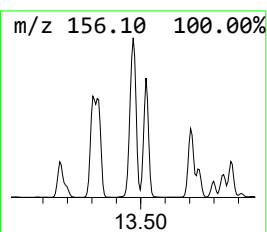
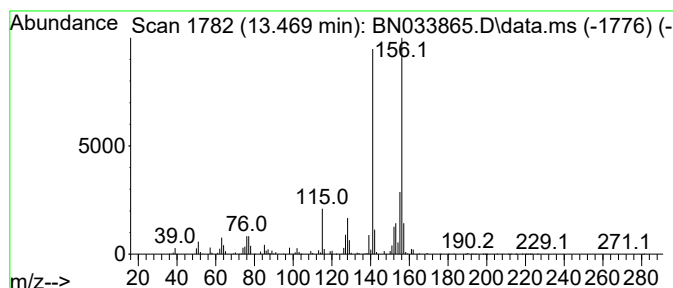
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 30 Naphthalene, 2,3-dimethyl- Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.469	7.95 ng/ul	8363060	Acenaphthene-d10	14.145

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
4			Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	97
5			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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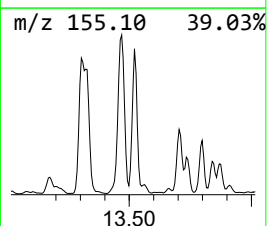
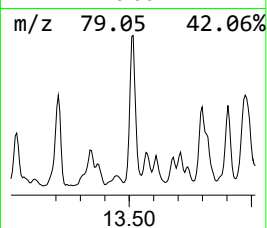
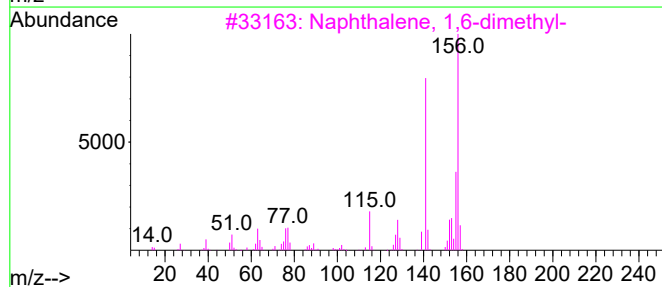
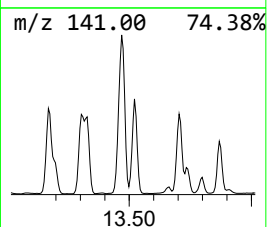
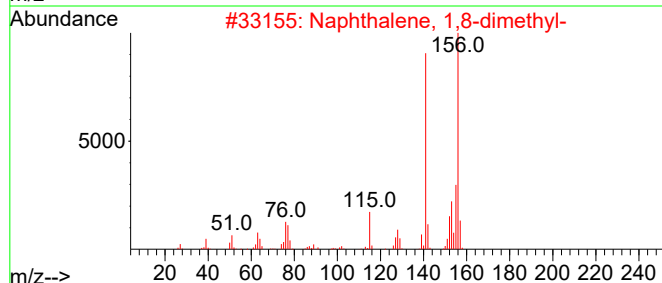
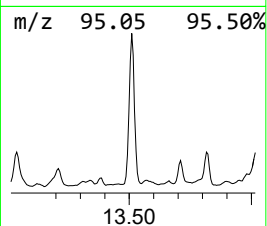
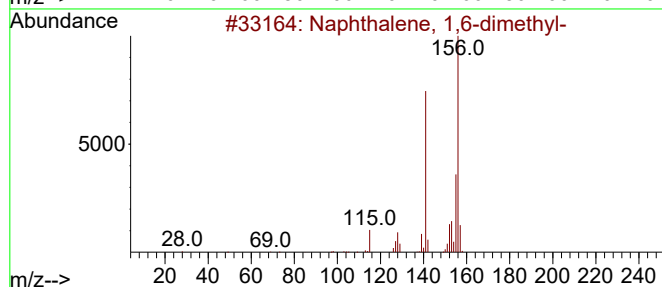
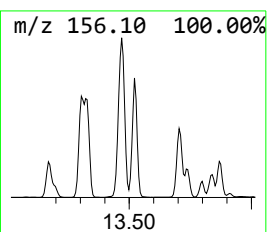
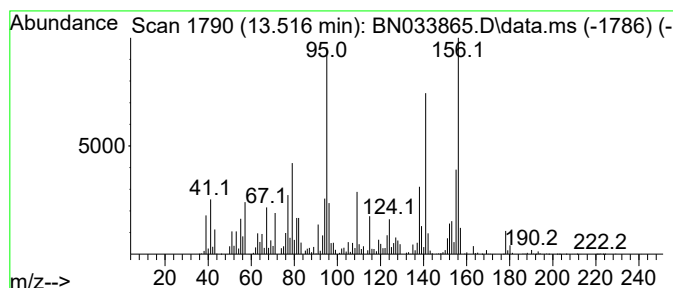
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 31 Naphthalene, 1,8-dimethyl- Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.516	10.15 ng/ul	10685900	Acenaphthene-d10	14.145

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	90
2			Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	83
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	83
4			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	83
5			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
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Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

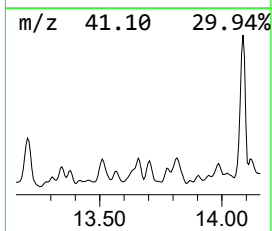
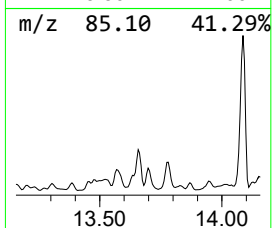
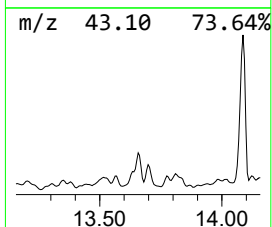
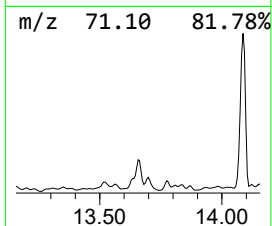
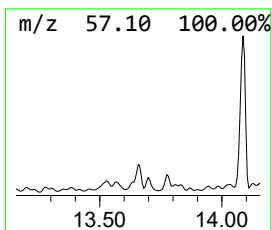
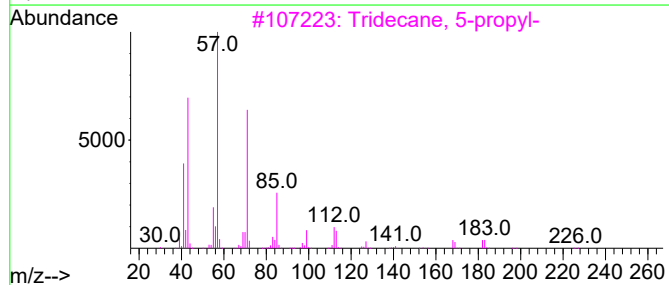
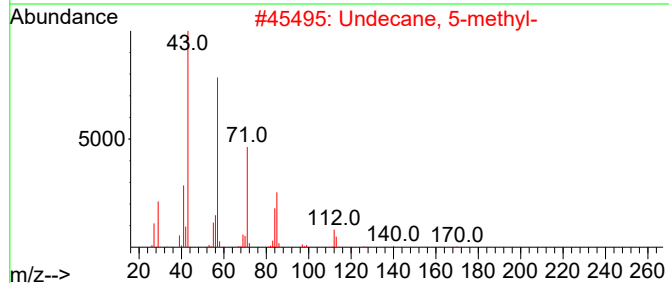
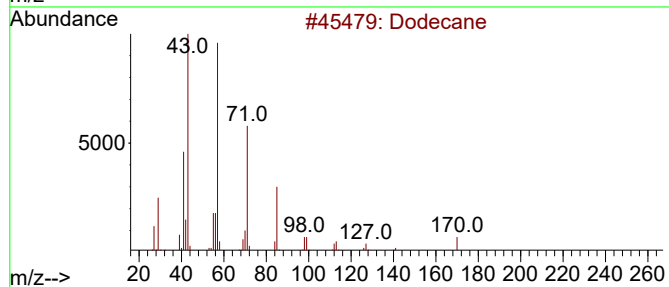
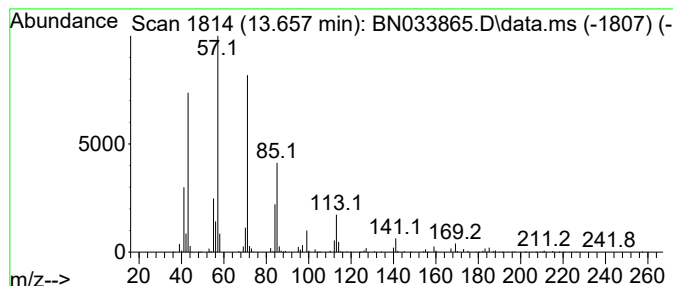
Instrument :
BNA_N
ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 60

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.657	4.26 ng/ul	4482250	Acenaphthene-d10	14.145	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane	170	C12H26	000112-40-3	74
2	Undecane, 5-methyl-	170	C12H26	001632-70-8	72
3	Tridecane, 5-propyl-	226	C16H34	055045-11-9	64
4	Dodecane	170	C12H26	000112-40-3	47
5	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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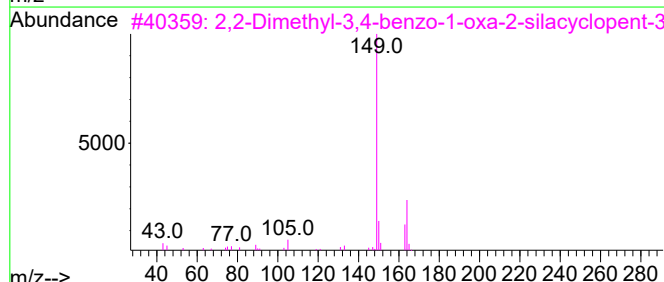
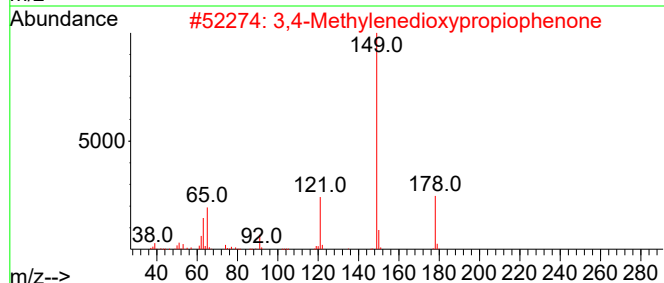
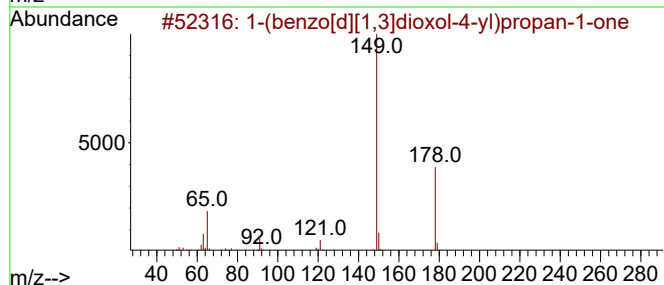
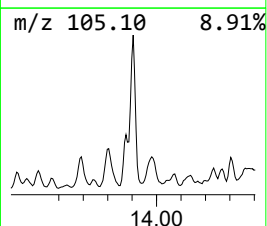
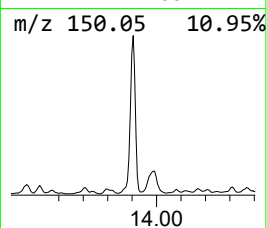
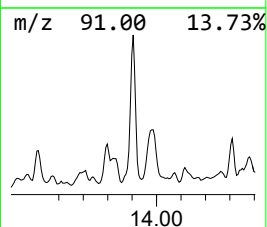
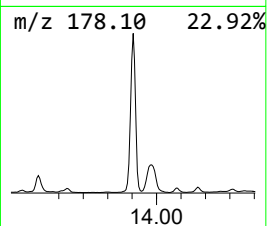
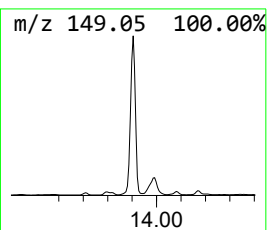
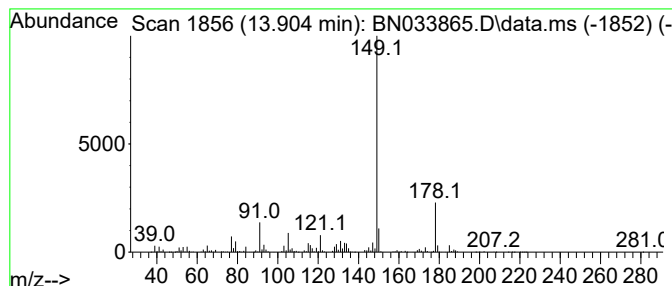
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 34 1-(benzo[d][1,3]dioxol-4-yl)... Concentration Rank 51

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.904	7.43 ng/ul	7816270	Acenaphthene-d10	14.145

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-(benzo[d][1,3]dioxol-4-yl)prop...	178	C10H10O3	1000456-65-5	64
2		3,4-Methylenedioxypropiphenone	178	C10H10O3	028281-49-4	59
3		2,2-Dimethyl-3,4-benzo-1-oxa-2-s...	164	C9H12OSi	1000424-28-3	59
4		3,4-Methylenedioxypropiphenone	178	C10H10O3	028281-49-4	59
5		Silane, 1,3-heptadiynyltrimethyl-	164	C10H16Si	019024-47-6	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

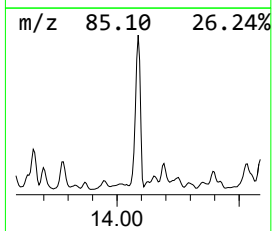
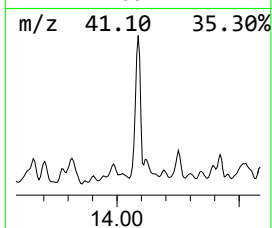
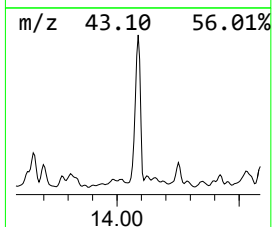
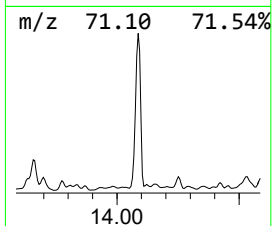
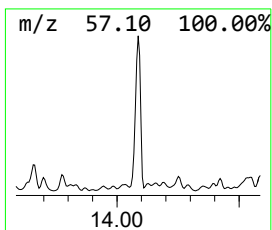
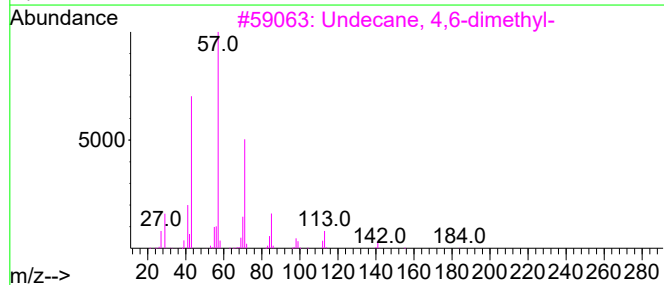
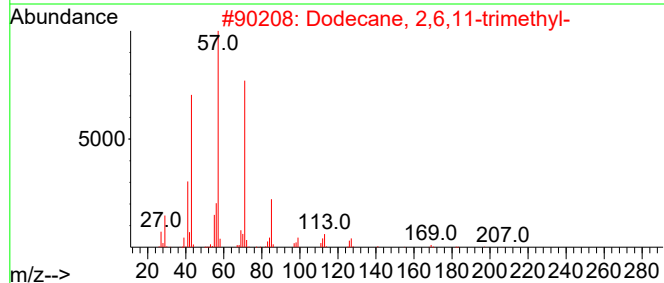
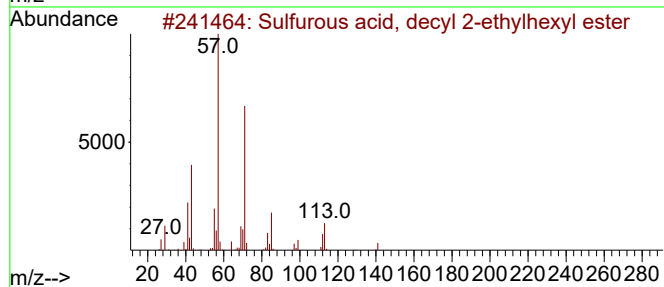
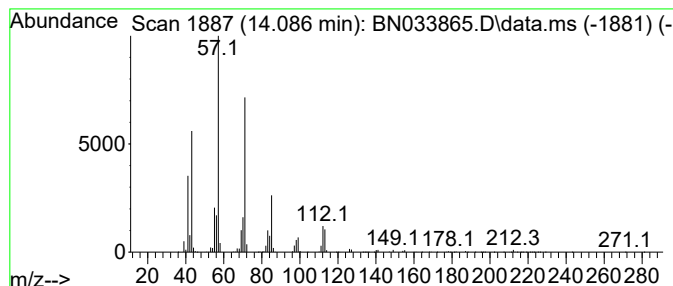
Instrument :
BNA_N
ClientSampleId :
DDCC1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 36 Sulfurous acid, decyl 2-eth... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.086	20.00 ng/ul	21052100	Acenaphthene-d10	14.145	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Sulfurous acid, decyl 2-ethylhex...	334	C18H38O3S	1000309-19-3	80
2	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	80
3	Undecane, 4,6-dimethyl-	184	C13H28	017312-82-2	72
4	Pentadecane, 7-methyl-	226	C16H34	006165-40-8	72
5	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
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Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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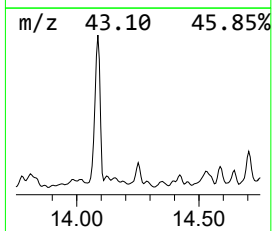
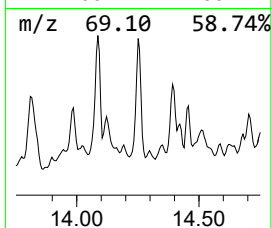
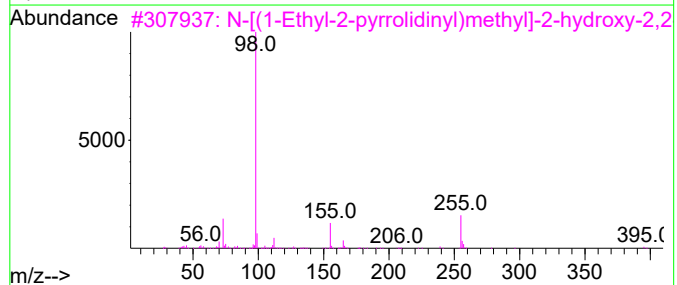
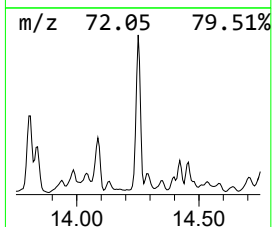
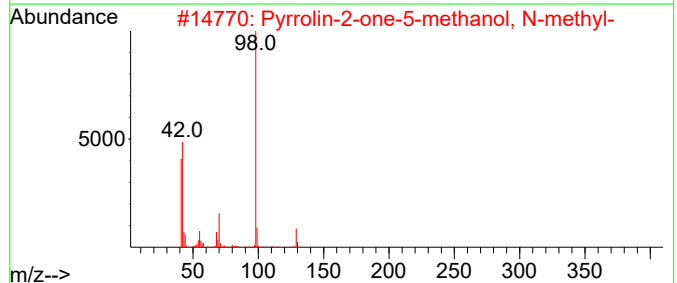
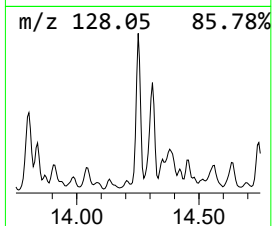
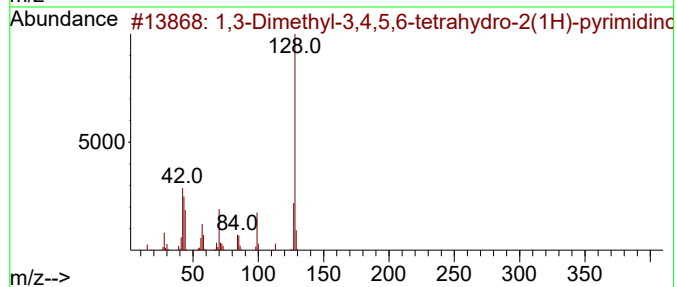
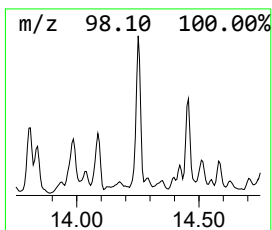
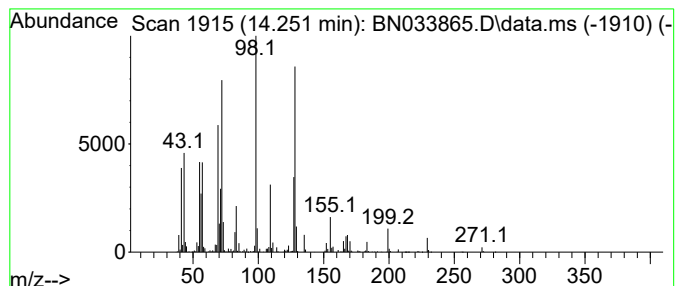
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 37 unknown-03 Concentration Rank 54

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.251	6.05 ng/ul	6363900	Acenaphthene-d10	14.145

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Dimethyl-3,4,5,6-tetrahydro-...	128	C6H12N2O	007226-23-5	22
2			Pyrrolin-2-one-5-methanol, N-met...	129	C6H11NO2	1000125-98-2	18
3			N-[(1-Ethyl-2-pyrrolidinyl)methy...	410	C24H34N2O2Si	1000505-32-8	14
4			2-(2-Chlorophenoxy)ethanol, isop...	214	C11H15ClO2	1000394-99-0	12
5			3-Undecanone	170	C11H22O	002216-87-7	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
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Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

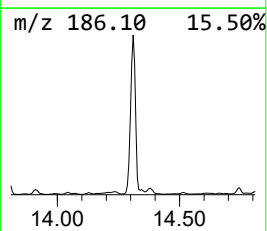
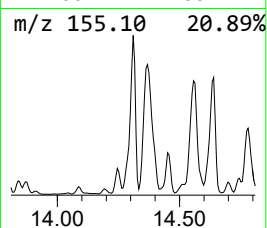
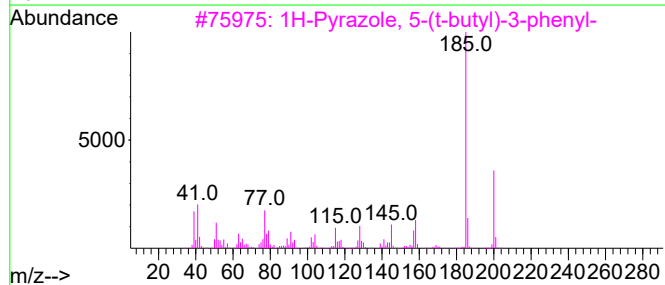
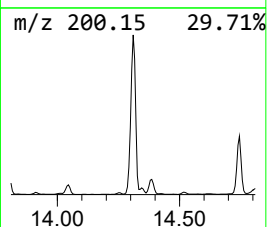
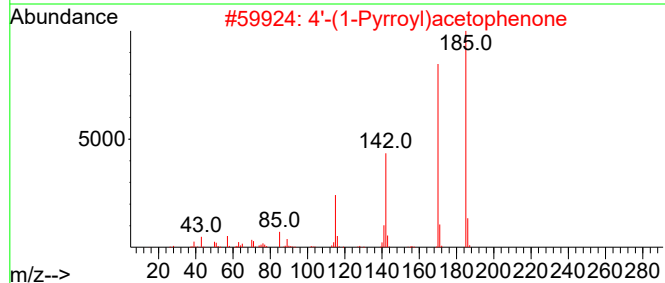
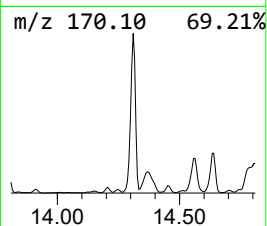
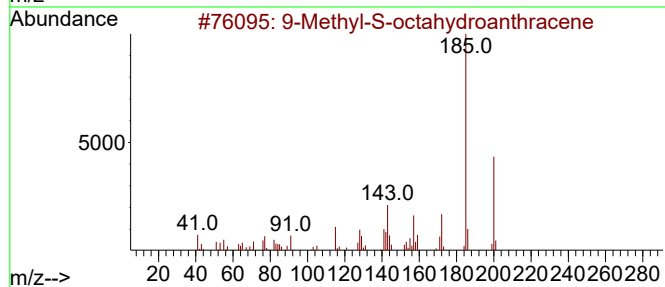
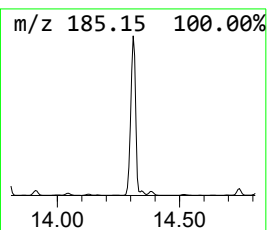
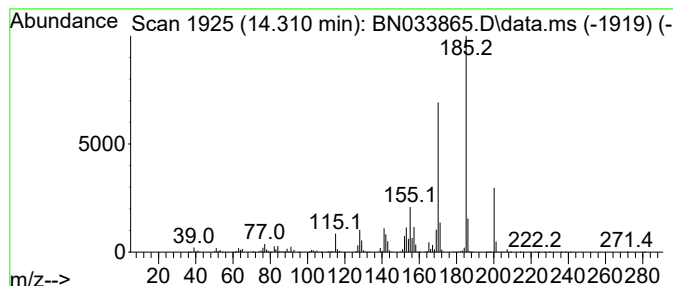
Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 38 9-Methyl-S-octahydroanthracene Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.310	15.74 ng/ul	16566100	Acenaphthene-d10	14.145	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Methyl-S-octahydroanthracene	200	C15H20	004948-51-0	60
2	4'-(1-Pyrroyl)acetophenone	185	C12H11NO	022106-37-2	58
3	1H-Pyrazole, 5-(t-butyl)-3-phenyl-	200	C13H16N2	1000261-34-8	55
4	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	51
5	Methyl 2-diethylamino-3-methyl-b...	185	C10H19NO2	075122-81-5	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
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Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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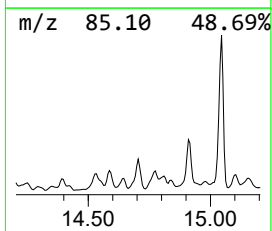
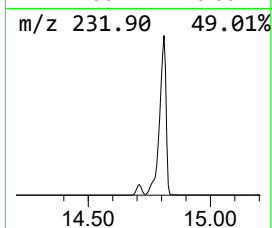
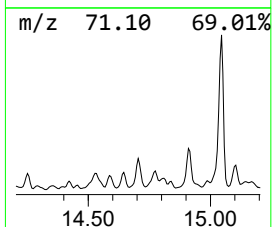
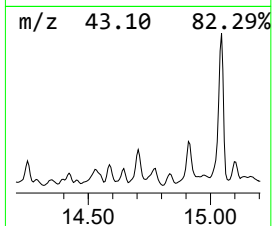
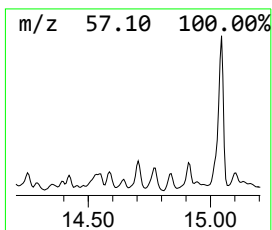
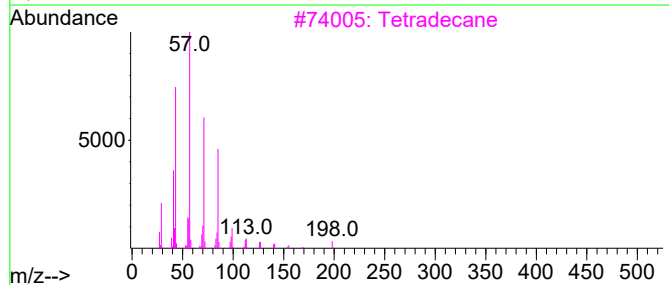
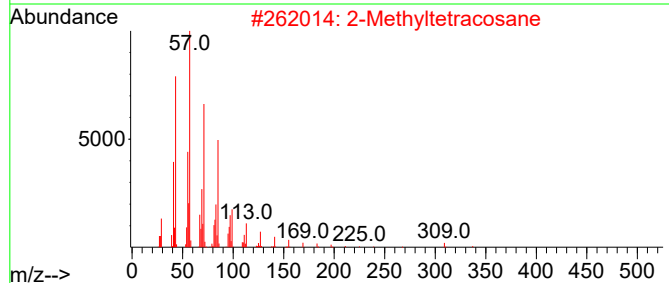
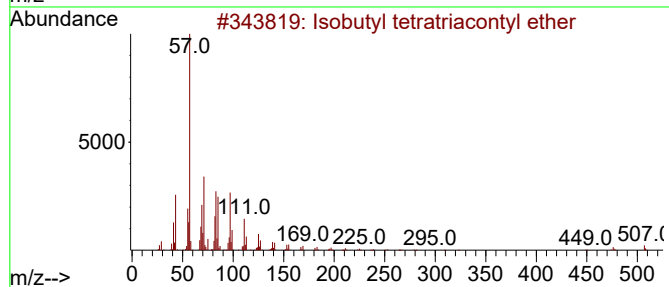
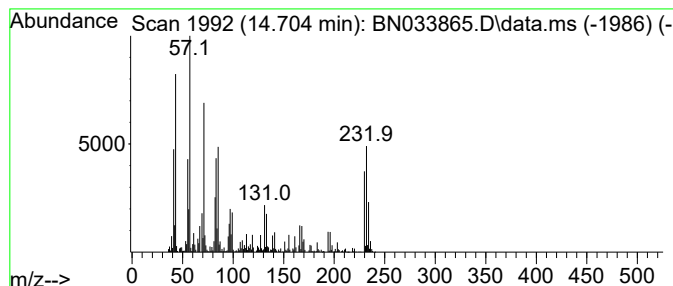
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 40 unknown-04 Concentration Rank 47

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.704	9.03 ng/ul	9501650	Acenaphthene-d10	14.145

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isobutyl tetratriacontyl ether	551	C38H78O	1000406-33-7	38
2			2-Methyltetracosane	352	C25H52	001560-78-7	38
3			Tetradecane	198	C14H30	000629-59-4	35
4			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	35
5			Dodecane, 2-methyl-	184	C13H28	001560-97-0	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

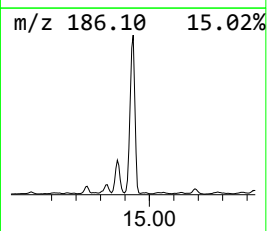
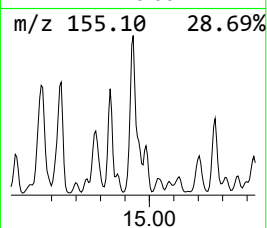
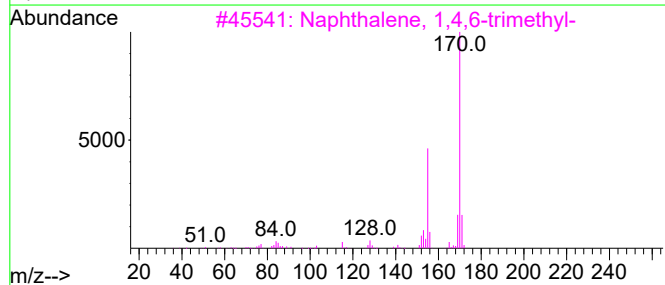
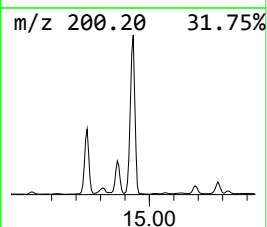
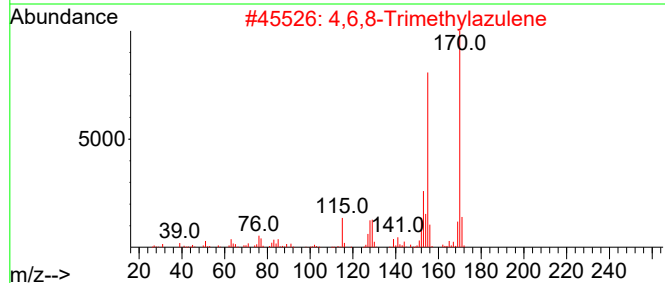
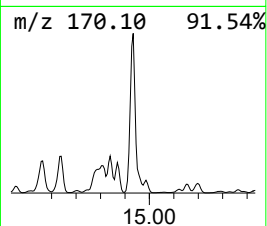
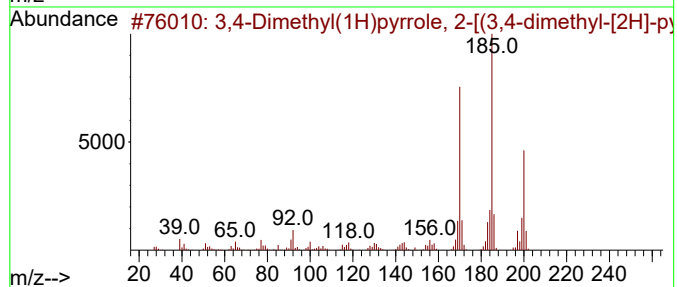
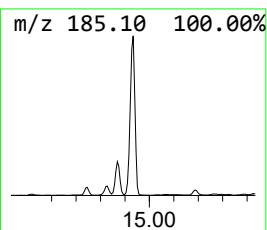
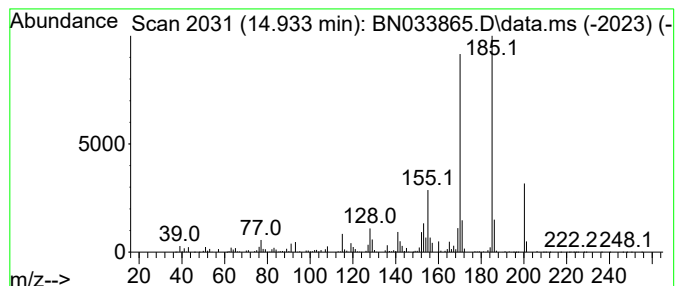
Instrument :
BNA_N
ClientSampleId :
DDCC1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 42 3,4-Dimethyl(1H)pyrrole, 2-... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.933	21.73 ng/ul	22868400	Acenaphthene-d10	14.145	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,4-Dimethyl(1H)pyrrole, 2-[(3,4...	200	C13H16N2	065539-79-9	72
2	4,6,8-Trimethylazulene	170	C13H14	000941-81-1	55
3	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	50
4	4,6,8-Trimethylazulene	170	C13H14	000941-81-1	50
5	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

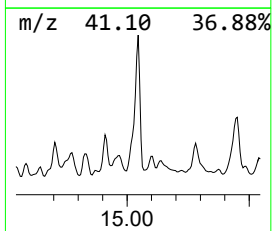
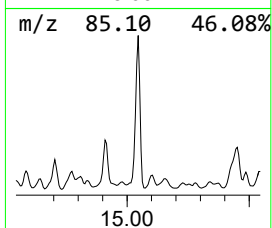
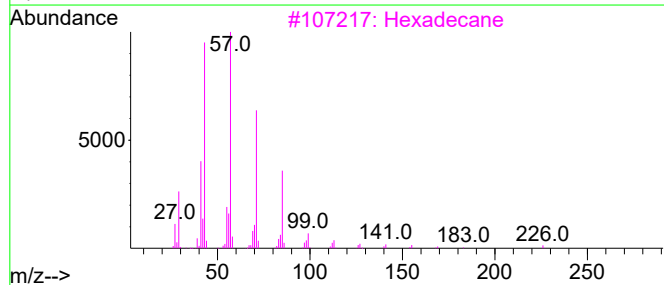
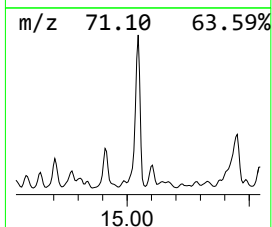
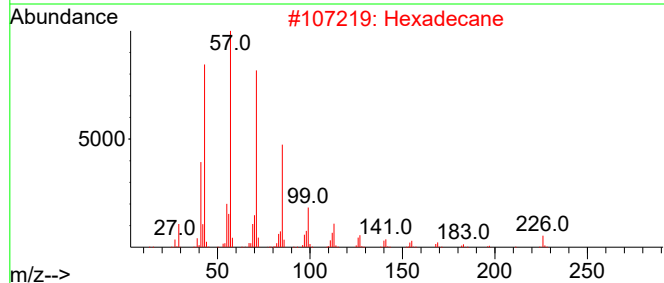
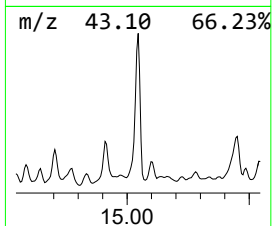
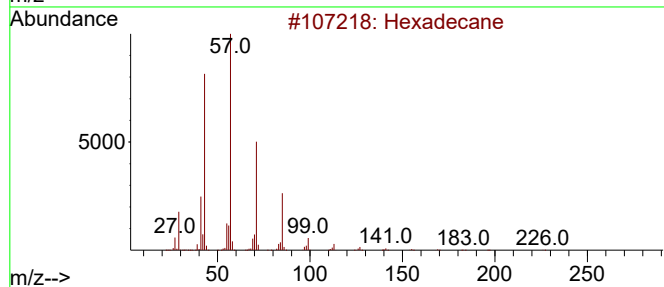
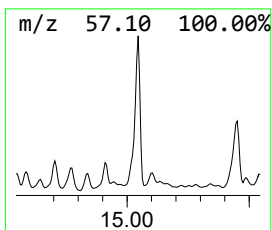
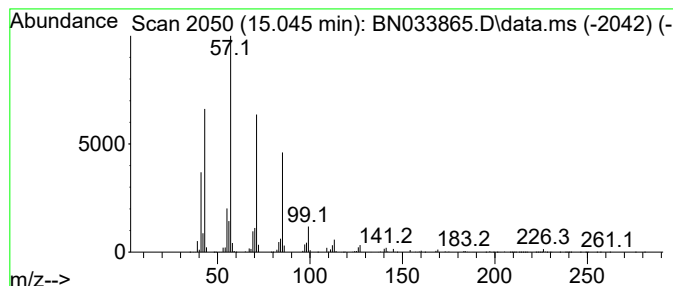
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 43 (DEL) Alkane: Straight-Chai... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.045	19.97 ng/ul	21025100	Acenaphthene-d10		14.145	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	96
2	Hexadecane		226	C16H34	000544-76-3	95
3	Hexadecane		226	C16H34	000544-76-3	94
4	Pentadecane		212	C15H32	000629-62-9	90
5	Nonadecane, 9-methyl-		282	C20H42	013287-24-6	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

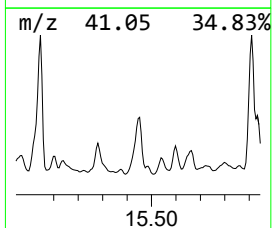
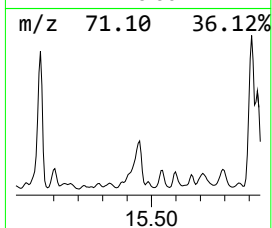
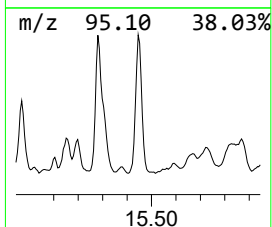
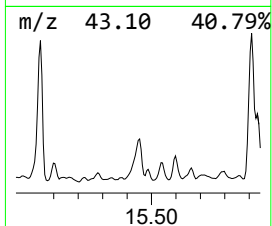
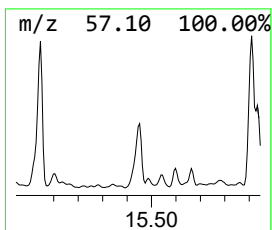
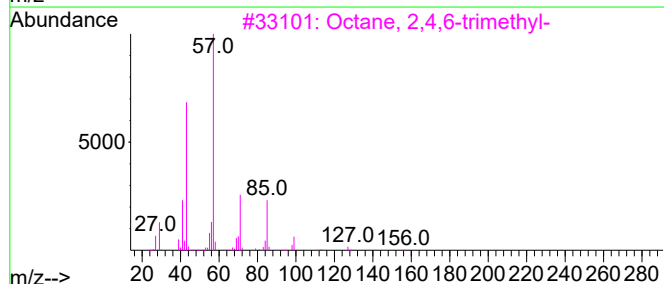
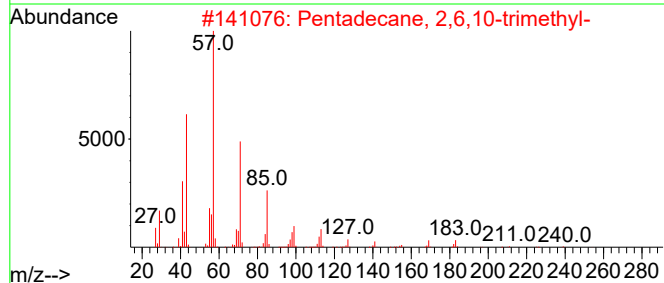
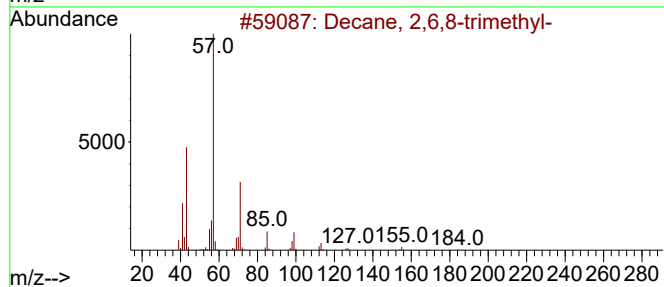
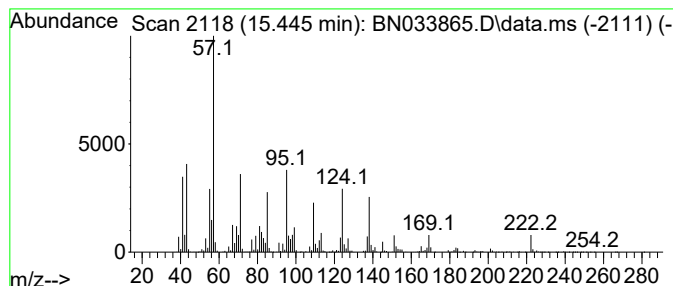
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 44 (DEL) Alkane: Straight-Chai... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD			R.T.
15.445	16.87 ng/ul	17757400	Acenaphthene-d10			14.145
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Decane, 2,6,8-trimethyl-		184	C13H28	062108-26-3	42
2	Pentadecane, 2,6,10-trimethyl-		254	C18H38	003892-00-0	35
3	Octane, 2,4,6-trimethyl-		156	C11H24	062016-37-9	35
4	Decane, 6-ethyl-2-methyl-		184	C13H28	062108-21-8	27
5	Pivalic acid, (9-borabicyclo[3.3...		222	C13H23BO2	088644-66-0	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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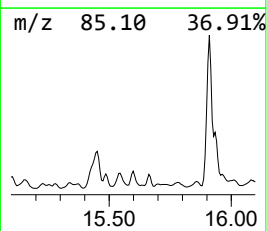
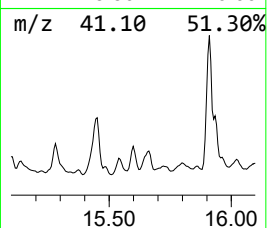
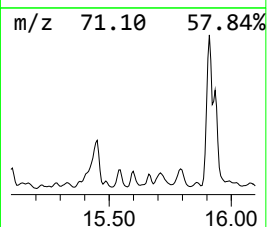
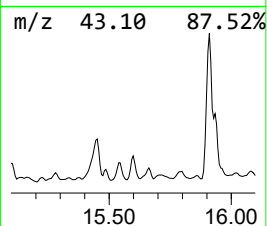
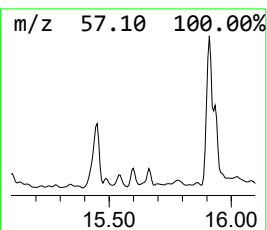
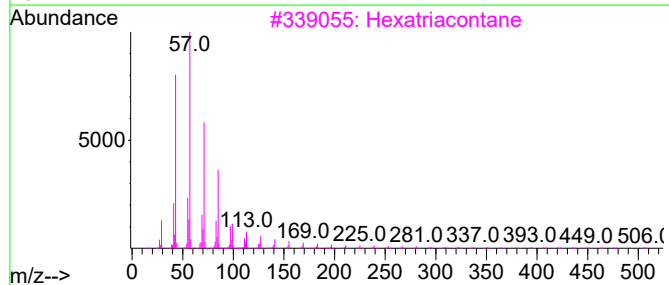
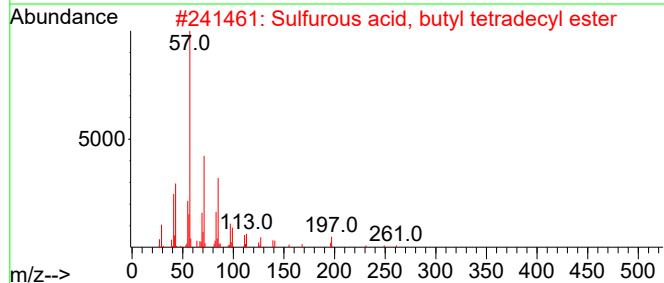
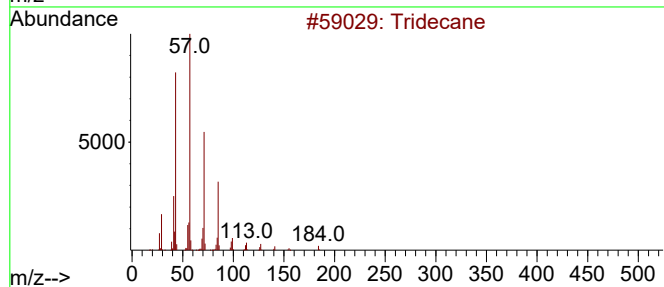
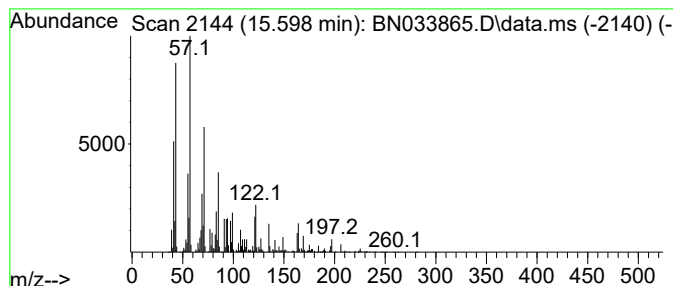
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 45 (DEL) Alkane: Straight-Chai... Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.598	8.47 ng/ul	4707280	Phenanthrene-d10	16.939

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	83
2			Sulfurous acid, butyl tetradecyl...	334	C18H38O3S	1010309-18-1	55
3			Hexatriacontane	507	C36H74	000630-06-8	46
4			Pentadecane	212	C15H32	000629-62-9	46
5			Sulfurous acid, dodecyl 2-propyl...	292	C15H32O3S	1000309-12-3	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DDCC1

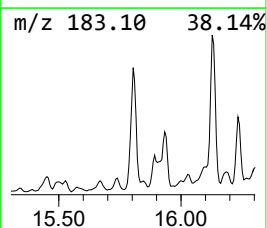
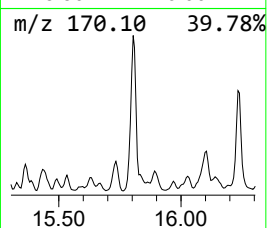
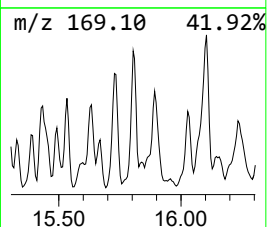
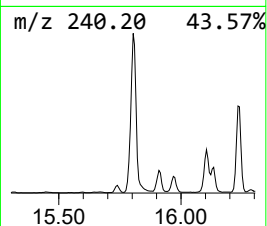
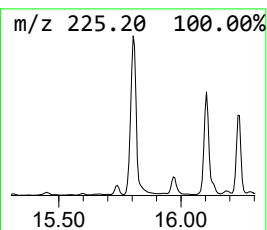
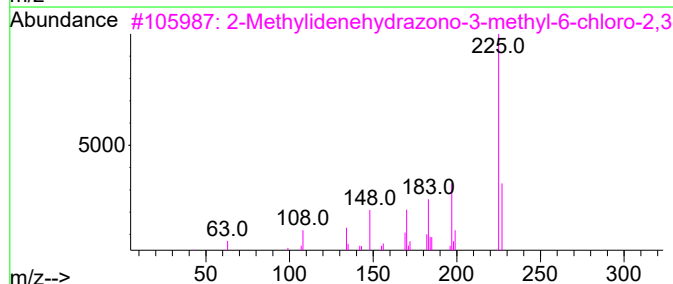
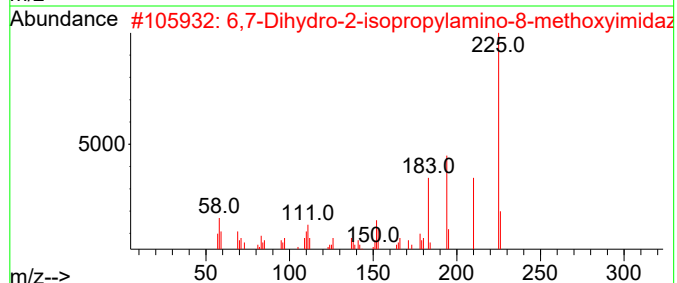
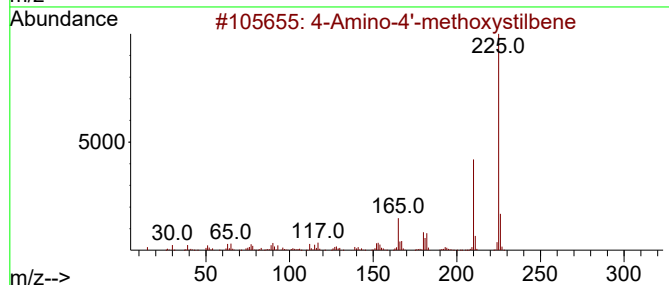
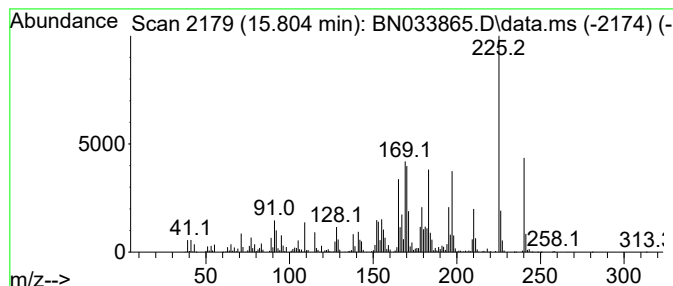
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 46 unknown-05 Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.804	16.77 ng/ul	9327280	Phenanthrene-d10	16.939

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Amino-4'-methoxystilbene	225	C15H15NO	007570-37-8	47
2			6,7-Dihydro-2-isopropylamino-8-m...	225	C9H15N5O2	1000101-98-2	46
3			2-Methylidenehydrazono-3-methyl-...	225	C9H8ClN3S	1000147-91-5	45
4			Naphthalene, 2,6-bis(1,1-dimethy...	240	C18H24	003905-64-4	38
5			(3-Isopropyl-5-methylamino-2,4,6...	225	C14H15N3	014204-00-3	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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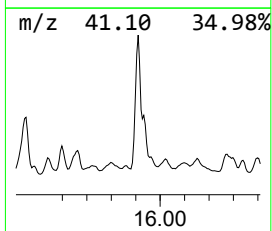
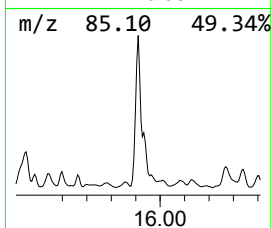
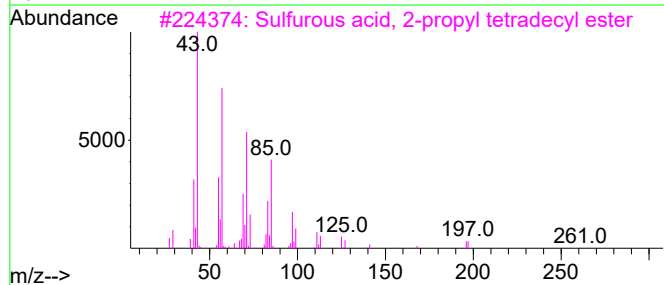
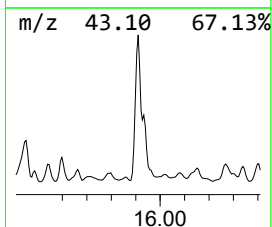
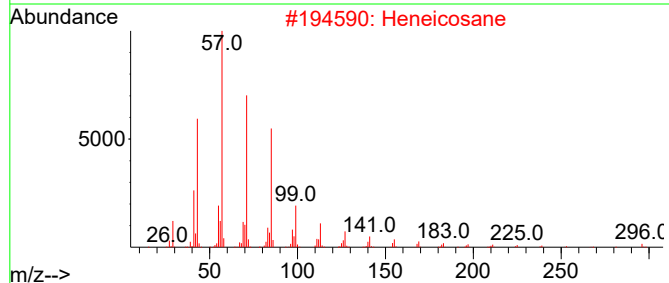
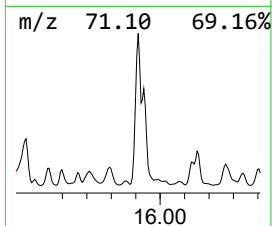
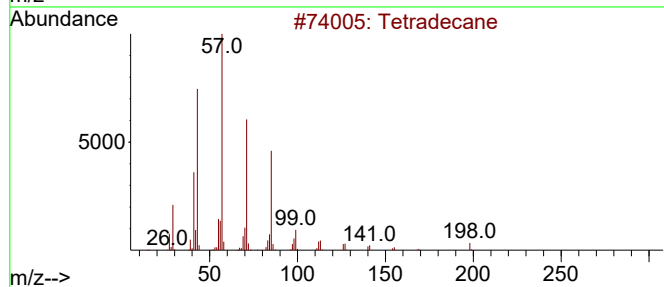
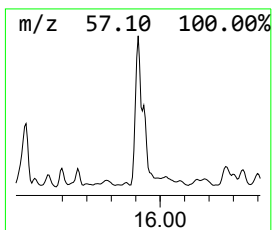
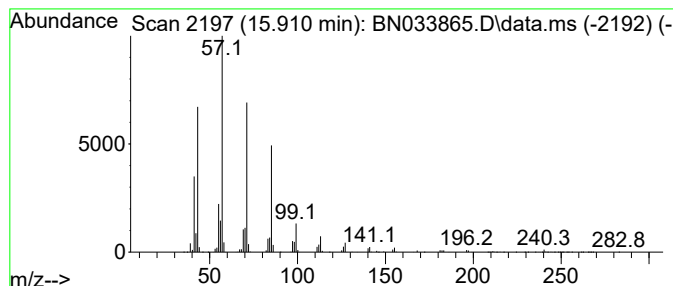
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 47 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.910	36.52 ng/ul	20305600	Phenanthrene-d10	16.939

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	91
2			Heneicosane	296	C21H44	000629-94-7	91
3			Sulfurous acid, 2-propyl tetrade...	320	C17H36O3S	1010309-12-5	90
4			Hexadecane	226	C16H34	000544-76-3	90
5			Tetradecane	198	C14H30	000629-59-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

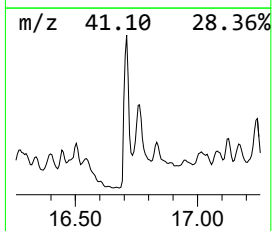
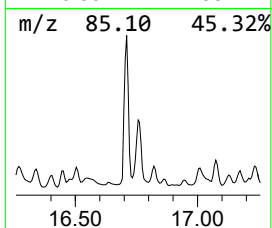
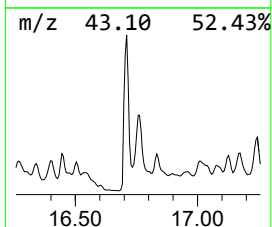
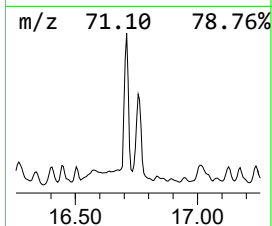
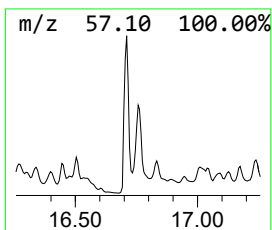
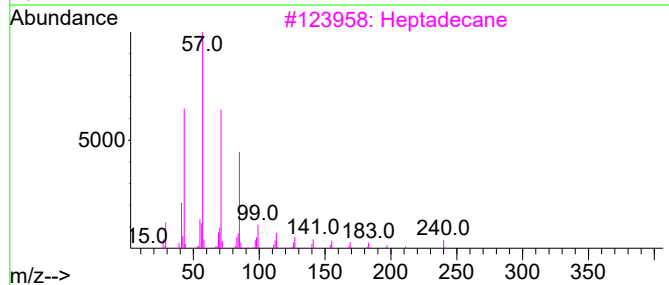
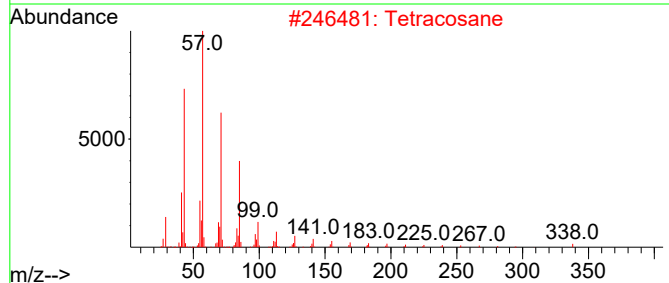
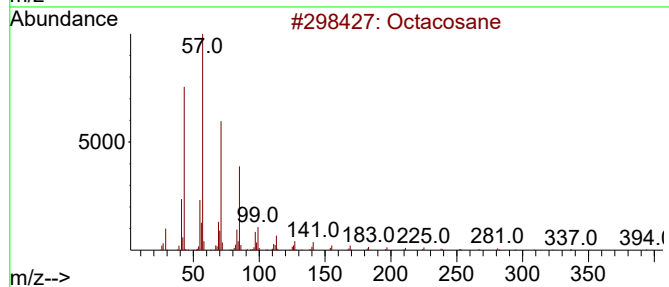
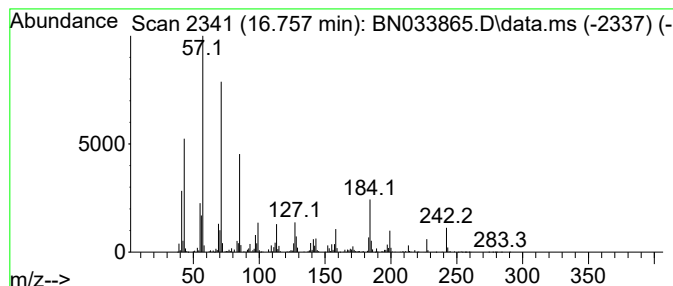
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ClientSampleId :
DDCC1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 48 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.757	28.32 ng/ul	15744800	Phenanthrene-d10		16.939	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Octacosane		394	C28H58	000630-02-4	68
2	Tetracosane		338	C24H50	000646-31-1	64
3	Heptadecane		240	C17H36	000629-78-7	64
4	Undecane, 2-methyl-		170	C12H26	007045-71-8	64
5	Tetradecane		198	C14H30	000629-59-4	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

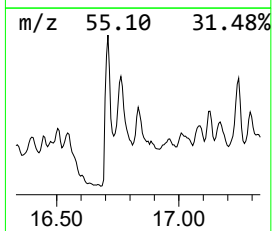
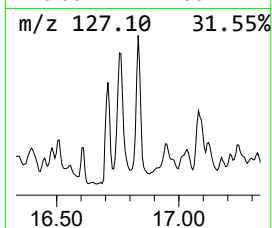
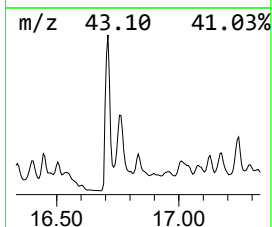
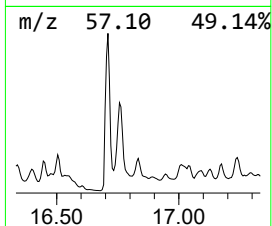
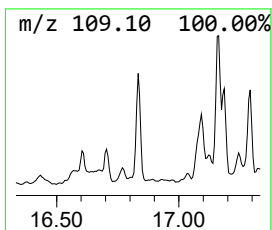
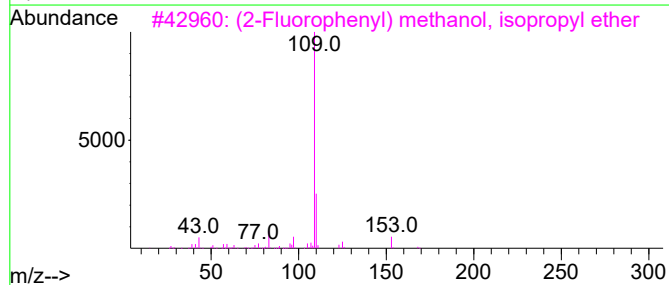
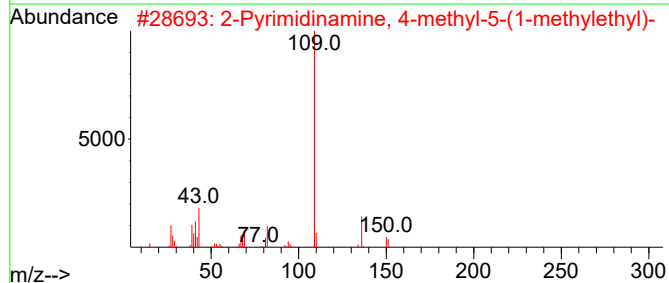
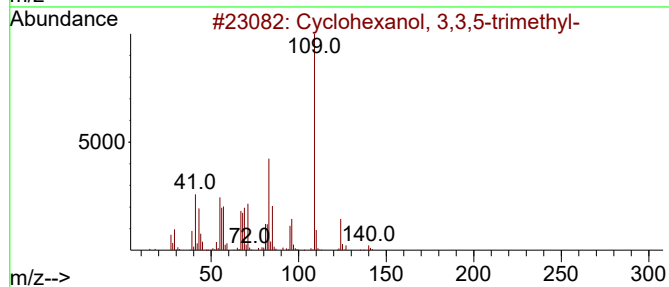
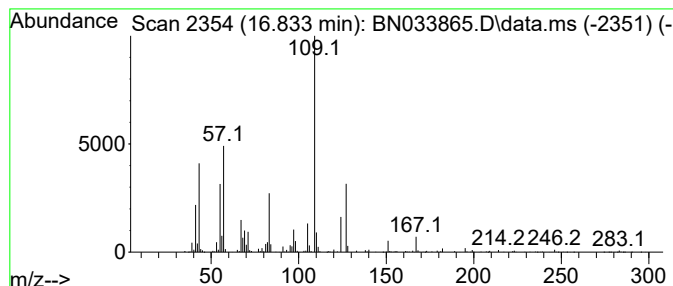
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 49 unknown-06 Concentration Rank 52

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.833	7.23 ng/ul	4020850	Phenanthrene-d10	16.939	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanol, 3,3,5-trimethyl-	142	C9H18O	000116-02-9	47
2	2-Pyrimidinamine, 4-methyl-5-(1-...	151	C8H13N3	1000460-76-8	43
3	(2-Fluorophenyl) methanol, isopr...	168	C10H13FO	1000374-56-3	43
4	2,5-Dimethylhex-5-en-3-yn-2-ol	124	C8H12O	1000302-74-9	43
5	2-Amino-4-methylpyrimidine	109	C5H7N3	000108-52-1	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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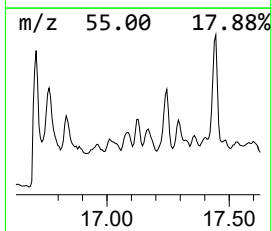
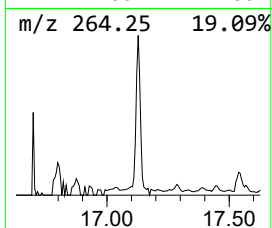
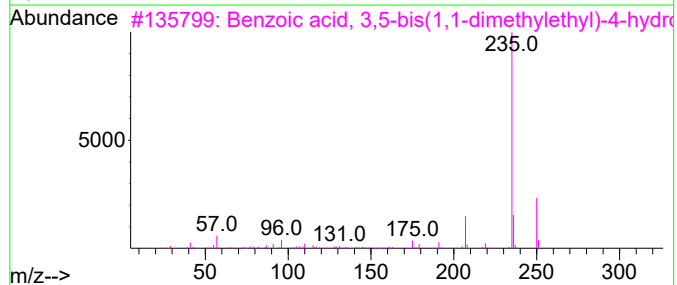
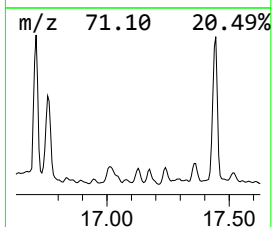
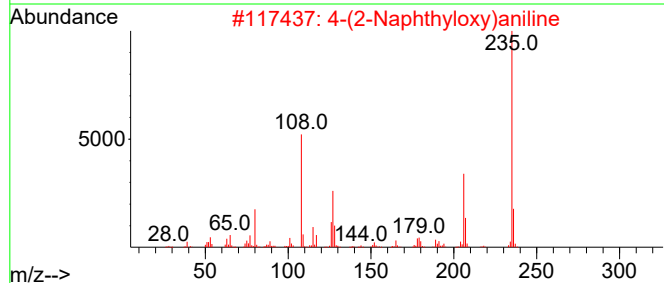
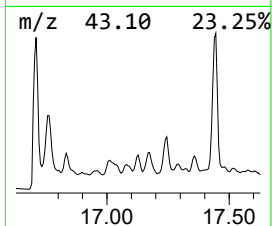
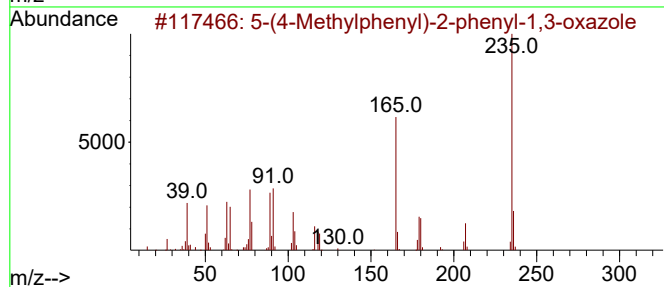
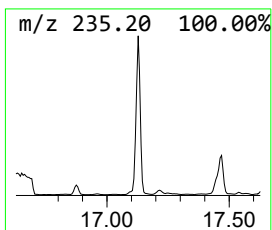
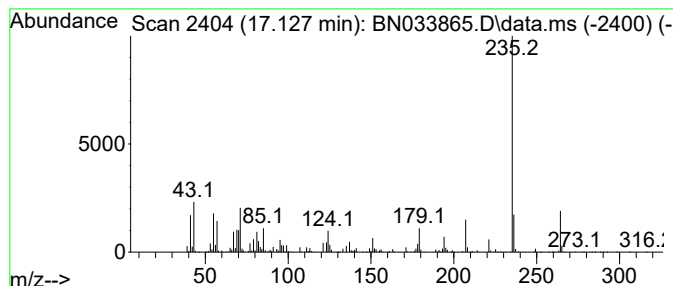
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 51 unknown-07 Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.127	9.30 ng/ul	5173180	Phenanthrene-d10	16.939

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-(4-Methylphenyl)-2-phenyl-1,3-...	235	C16H13NO	1000487-13-3	49		
2	4-(2-Naphthyloxy)aniline	235	C16H13NO	071311-83-6	47		
3	Benzoic acid, 3,5-bis(1,1-dimeth...	250	C15H22O3	001421-49-4	43		
4	Pyrazole, 5-amino-1,3-diphenyl-	235	C15H13N3	005356-71-8	43		
5	Pyrazole, 5-amino-1,3-diphenyl-	235	C15H13N3	005356-71-8	43		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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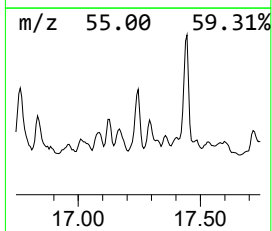
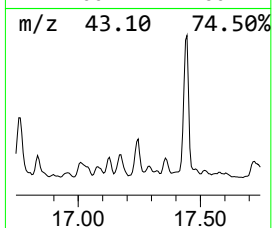
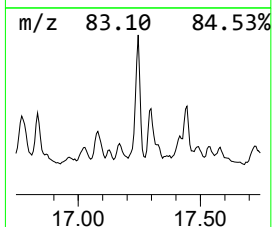
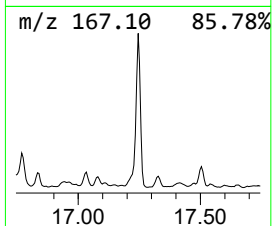
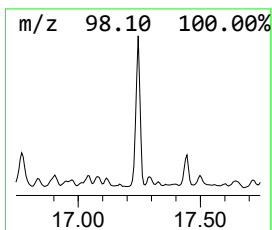
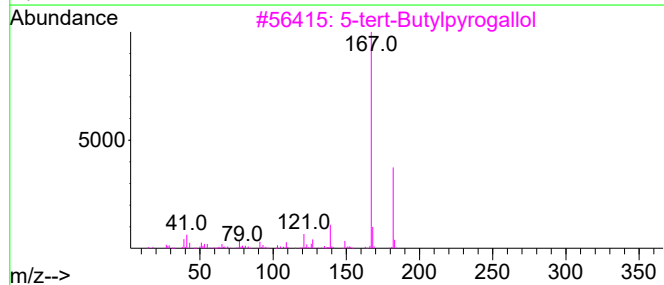
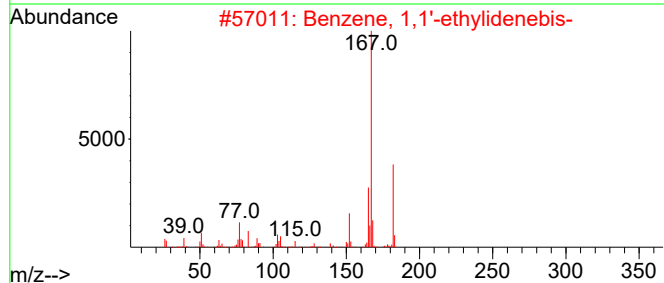
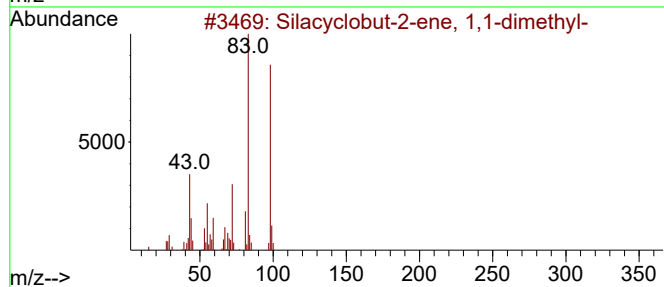
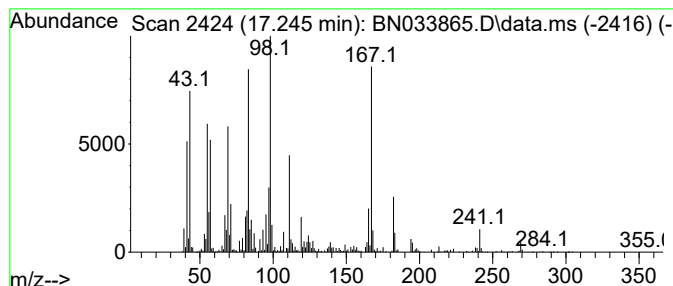
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 52 unknown-08 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.245	24.08 ng/ul	13387700	Phenanthrene-d10	16.939

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Silacyclobut-2-ene, 1,1-dimethyl-	98	C5H10Si	066222-35-3	38
2			Benzene, 1,1'-ethyldienebis-	182	C14H14	000612-00-0	25
3			5-tert-Butylpyrogallol	182	C10H14O3	020481-17-8	18
4			2-Cyclohexen-1-one, 3,5-dimethyl...	153	C9H15NO	056336-06-2	18
5			Cyclohexanone, 4-(1,1-dimethylpr...	168	C11H20O	016587-71-6	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
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Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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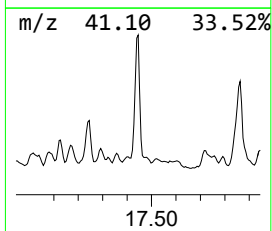
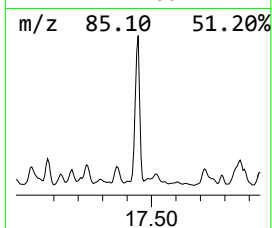
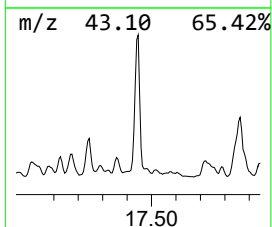
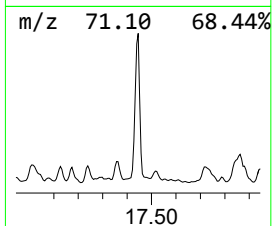
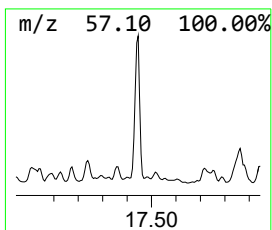
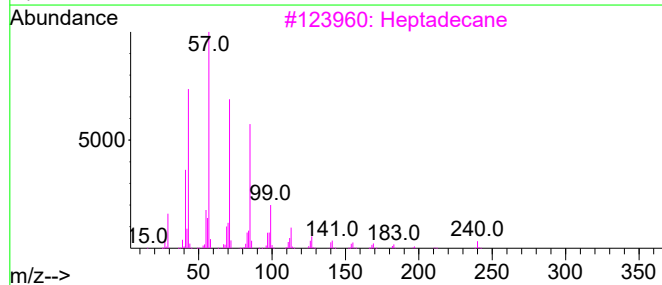
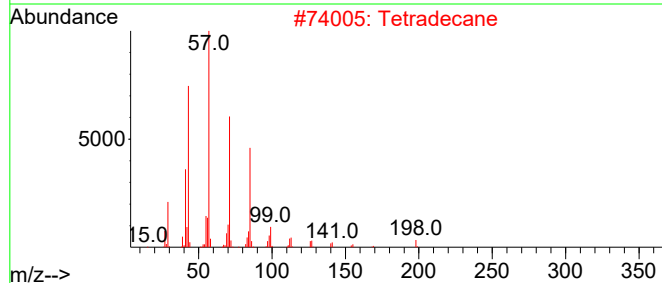
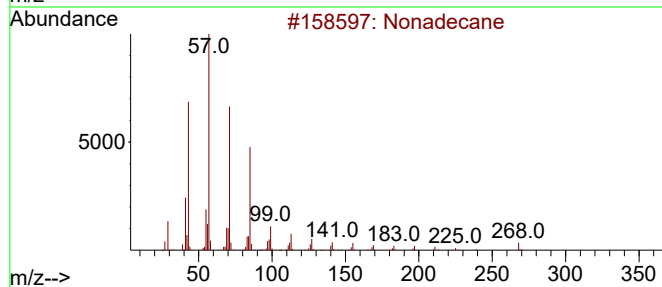
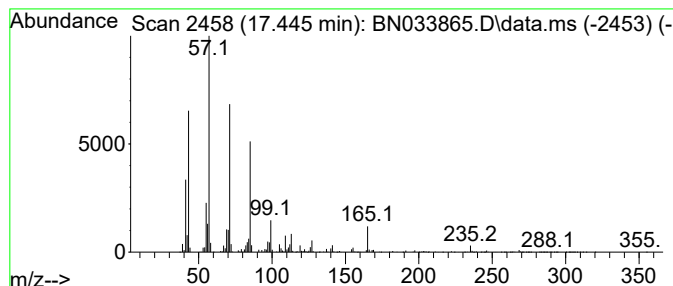
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 53 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.445	31.77 ng/ul	17668400	Phenanthrene-d10	16.939

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	94
2			Tetradecane	198	C14H30	000629-59-4	83
3			Heptadecane	240	C17H36	000629-78-7	83
4			Tetradecane	198	C14H30	000629-59-4	80
5			Hexadecane	226	C16H34	000544-76-3	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DDCC1

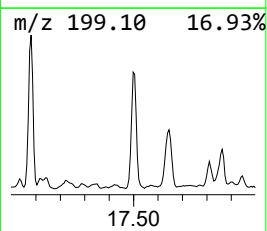
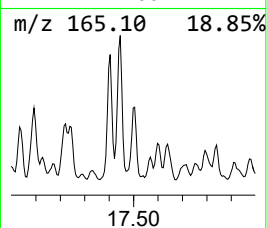
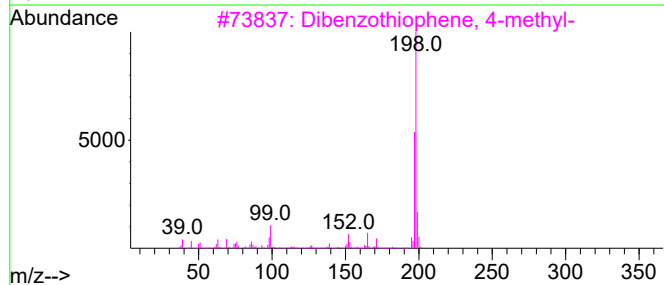
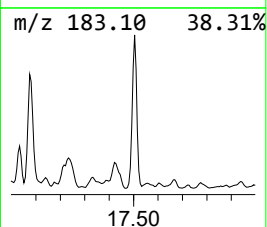
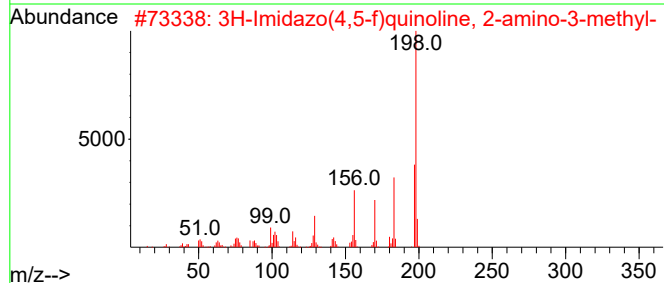
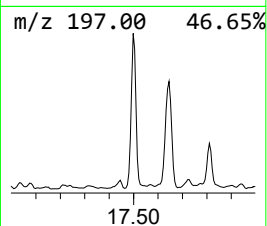
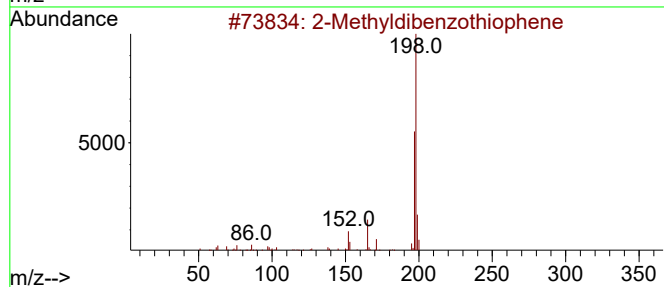
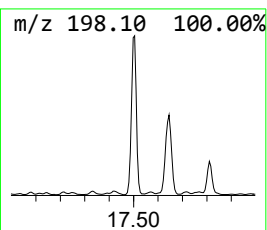
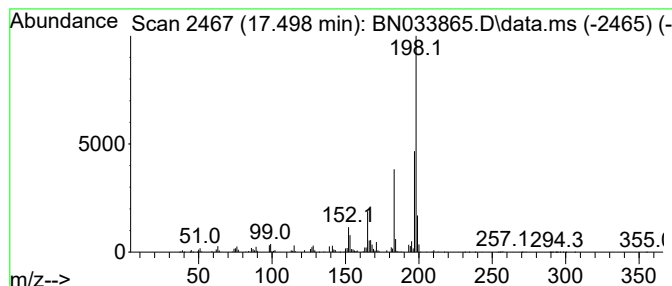
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 54 2-Methyldibenzothiophene Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.498	9.93 ng/ul	5523820	Phenanthrene-d10	16.939

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Methyldibenzothiophene	198	C13H10S	1000383-55-1	76
2			3H-Imidazo(4,5-f)quinoline, 2-am...	198	C11H10N4	076180-96-6	72
3			Dibenzothiophene, 4-methyl-	198	C13H10S	007372-88-5	62
4			Dibenzothiophene, 3-methyl-	198	C13H10S	016587-52-3	59
5			Benzene, 1-fluoro-4-(2-phenyleth...	198	C14H11F	000718-25-2	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
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Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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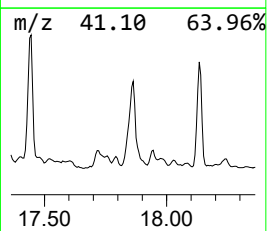
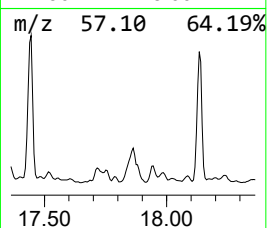
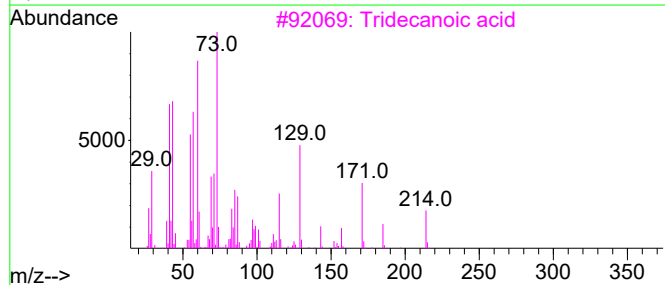
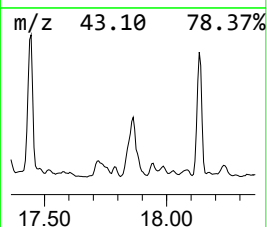
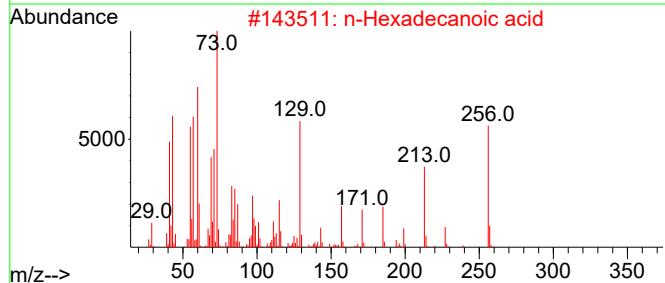
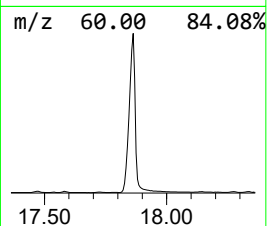
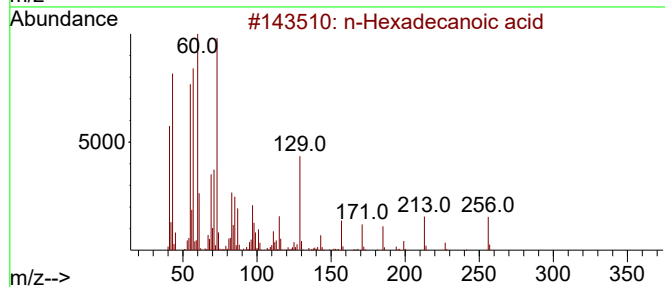
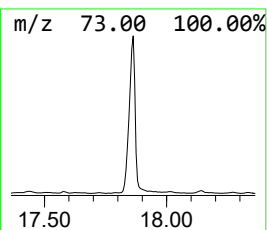
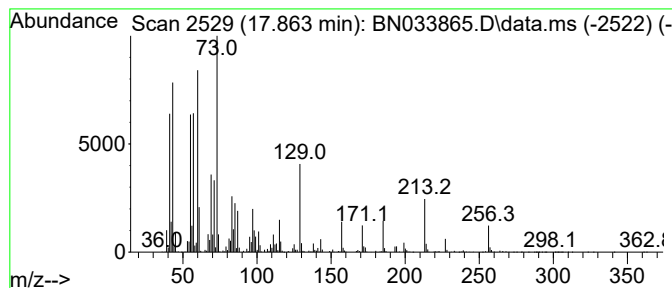
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 55 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.863	36.15 ng/ul	20102000	Phenanthrene-d10	16.939

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3			Tridecanoic acid	214	C13H26O2	000638-53-9	95
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91
5			Tridecanoic acid	214	C13H26O2	000638-53-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DDCC1

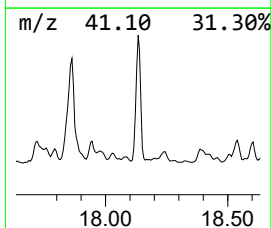
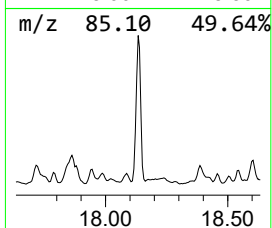
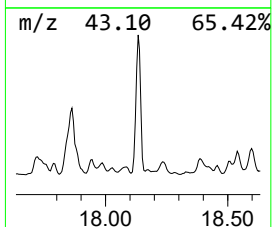
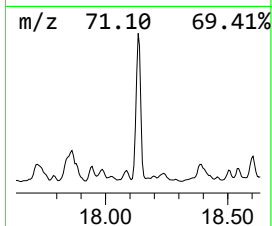
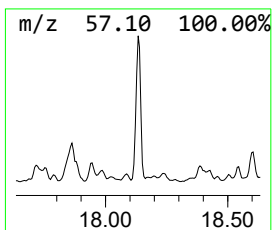
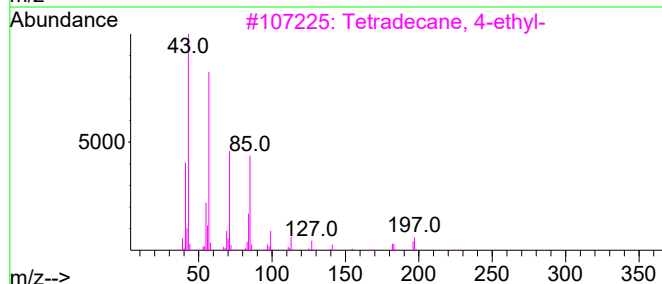
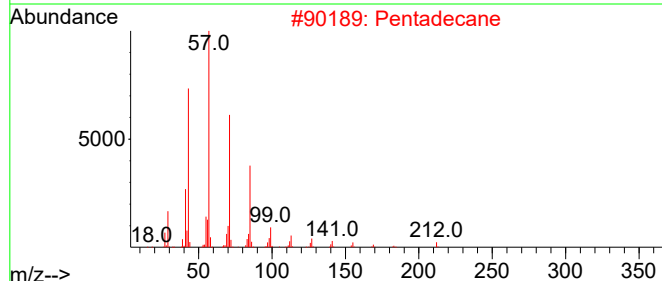
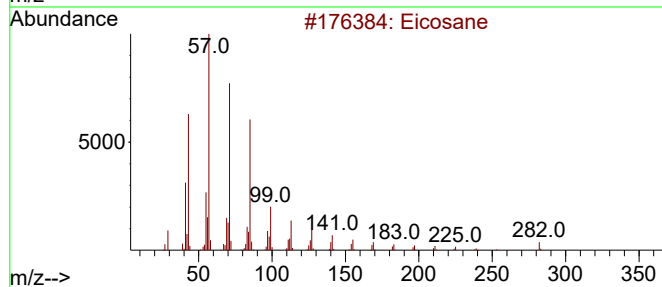
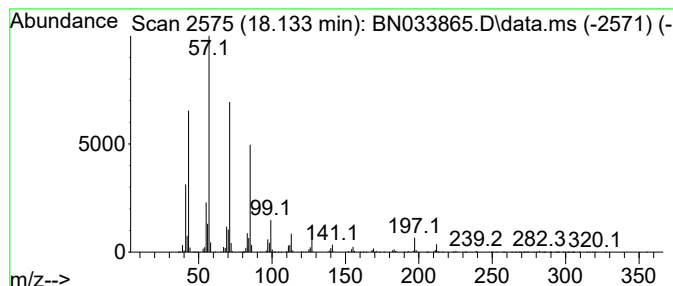
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 56 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.133	26.97 ng/ul	14994300	Phenanthrene-d10	16.939

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	97
2			Pentadecane	212	C15H32	000629-62-9	96
3			Tetradecane, 4-ethyl-	226	C16H34	055045-14-2	94
4			Pentadecane	212	C15H32	000629-62-9	92
5			Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DDCC1

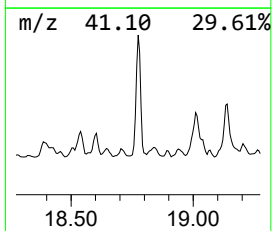
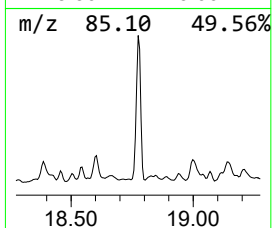
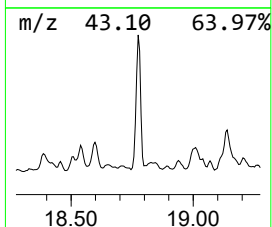
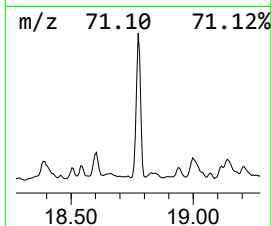
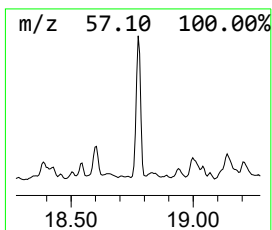
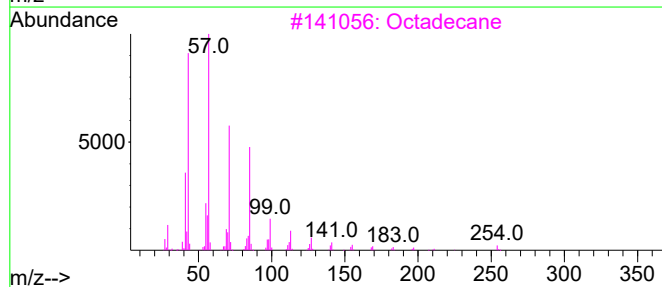
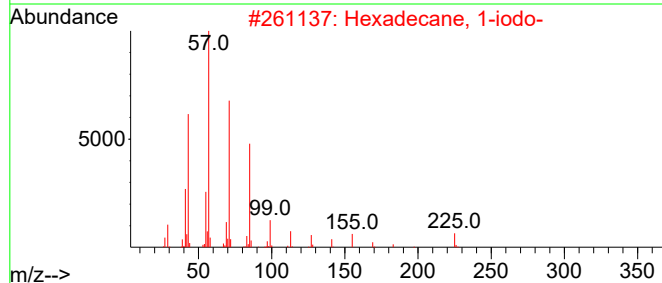
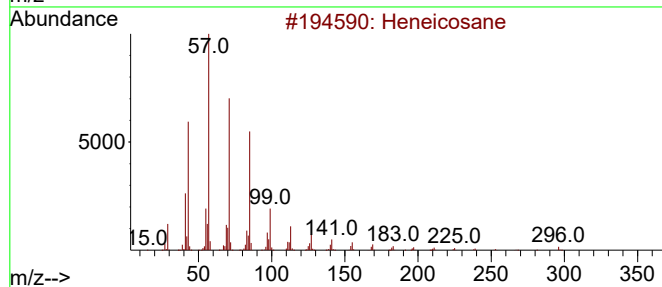
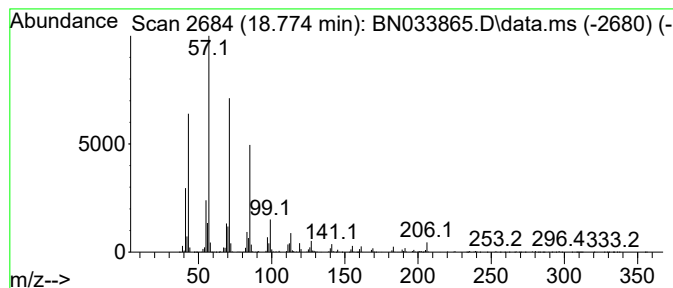
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 57 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.774	22.63 ng/ul	12580700	Phenanthrene-d10	16.939

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	90
2			Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	87
3			Octadecane	254	C18H38	000593-45-3	87
4			Nonadecane	268	C19H40	000629-92-5	87
5			Heptadecane	240	C17H36	000629-78-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DDCC1

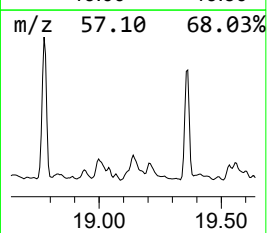
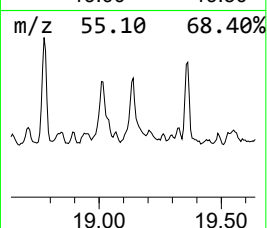
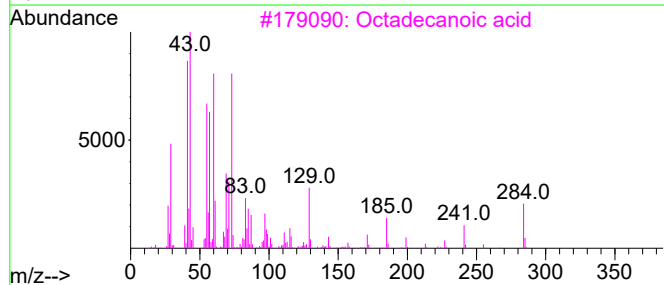
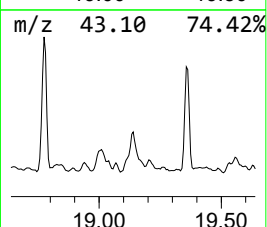
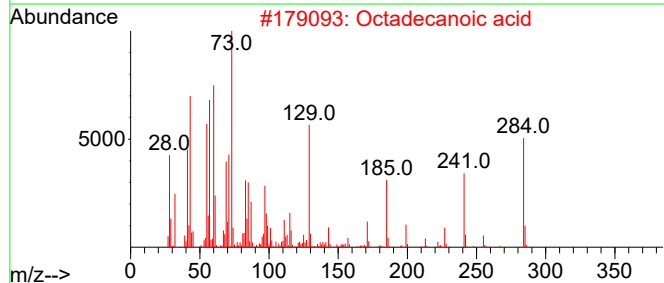
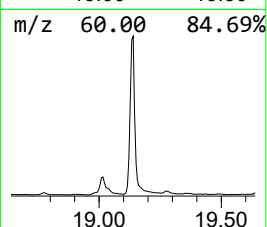
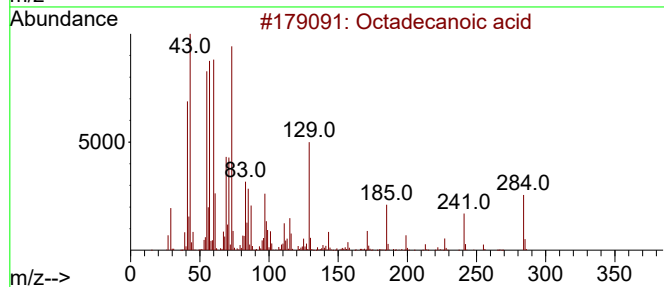
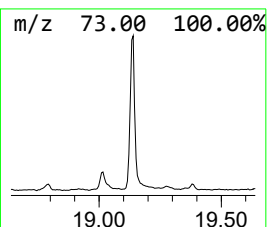
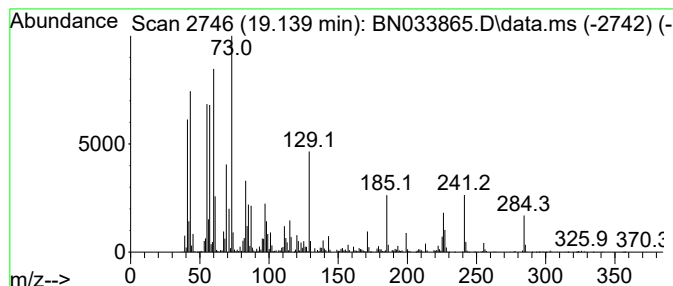
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 58 Octadecanoic acid Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.139	14.30 ng/ul	7505790	Chrysene-d12	21.139

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	96
2			Octadecanoic acid	284	C18H36O2	000057-11-4	95
3			Octadecanoic acid	284	C18H36O2	000057-11-4	94
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	93
5			Octadecanoic acid	284	C18H36O2	000057-11-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

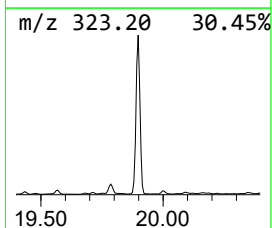
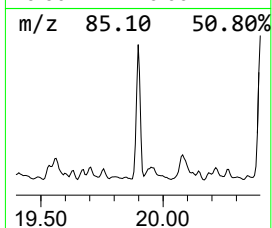
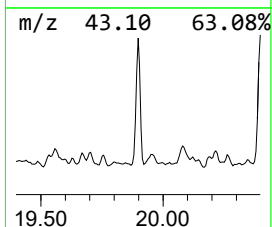
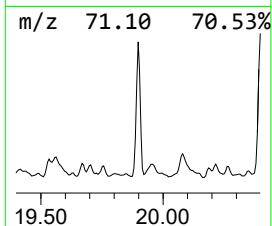
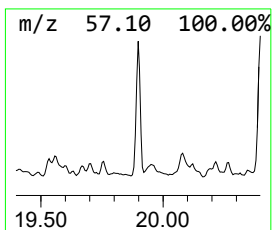
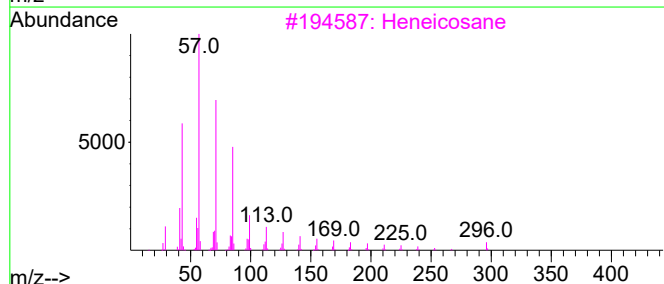
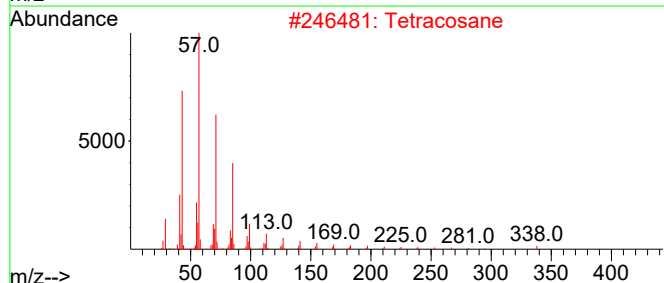
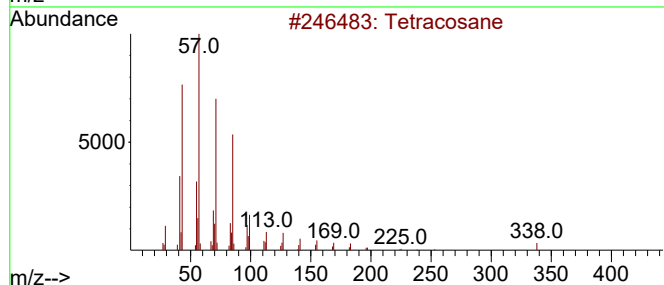
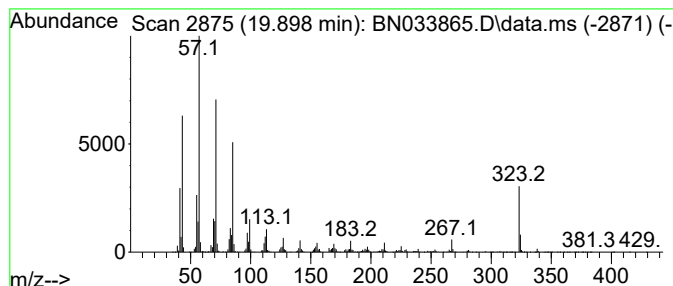
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 59 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.898	23.58 ng/ul	12377800	Chrysene-d12	21.139	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	98
2	Tetracosane	338	C24H50	000646-31-1	96
3	Heneicosane	296	C21H44	000629-94-7	95
4	Nonadecane	268	C19H40	000629-92-5	95
5	Nonadecane	268	C19H40	000629-92-5	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
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Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

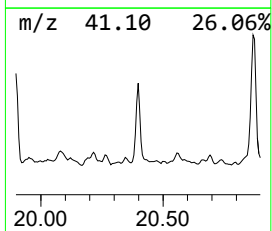
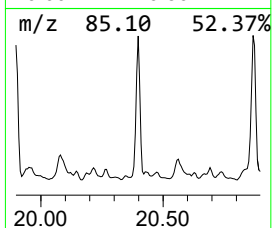
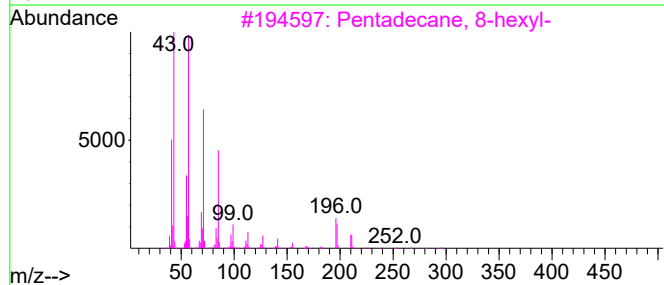
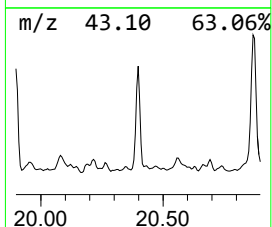
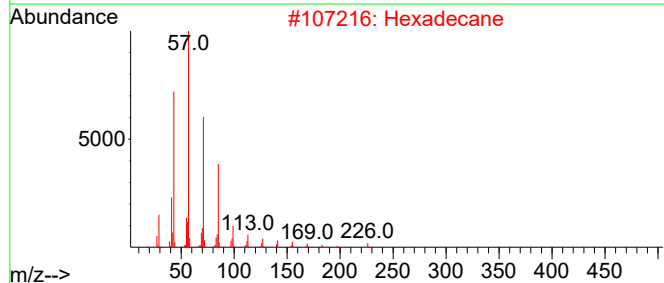
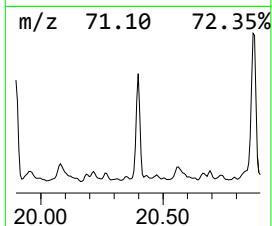
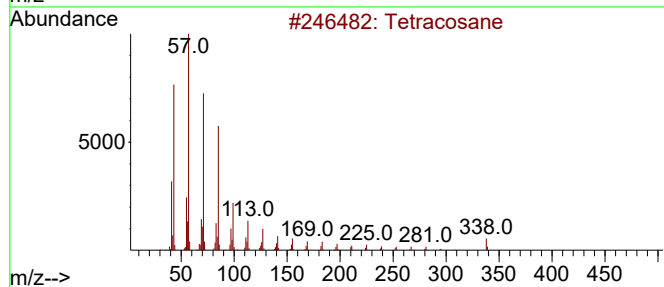
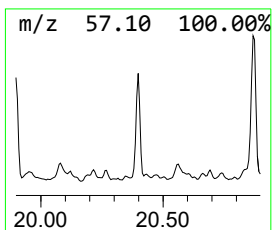
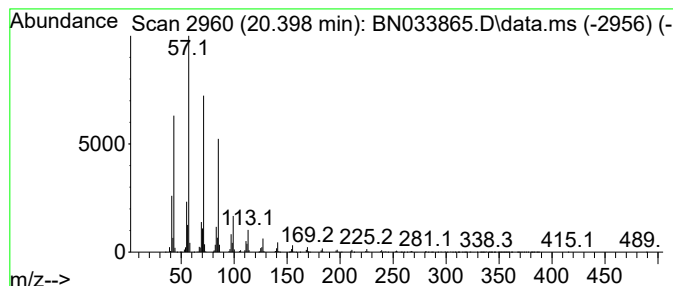
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 60 (DEL) Alkane: Straight-Chai... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.398	18.71 ng/ul	9822840	Chrysene-d12	21.139	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	97
2	Hexadecane	226	C16H34	000544-76-3	95
3	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	93
4	Nonadecane	268	C19H40	000629-92-5	93
5	Eicosane, 10-methyl-	296	C21H44	054833-23-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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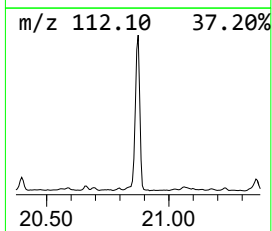
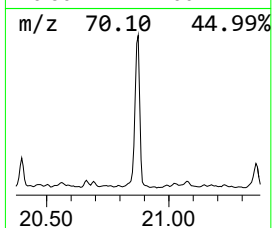
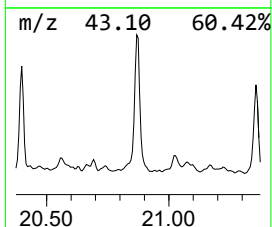
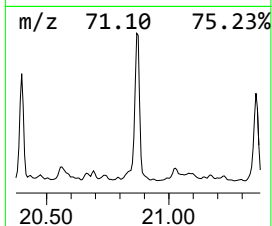
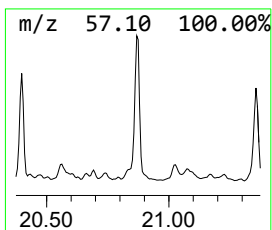
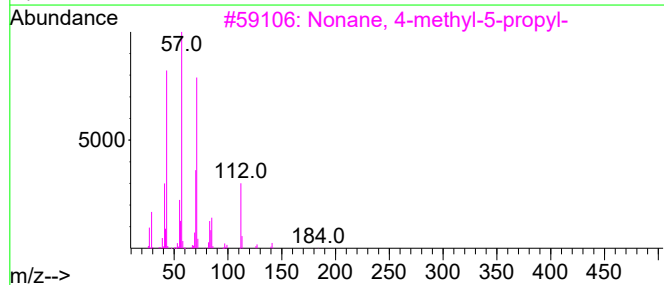
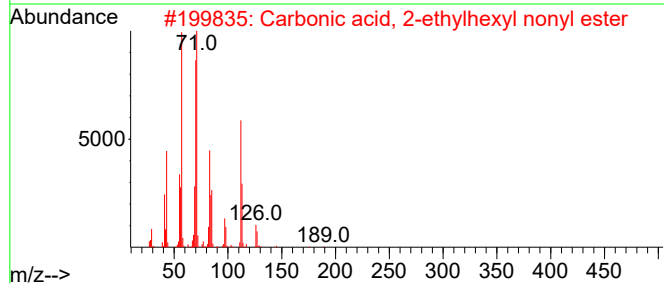
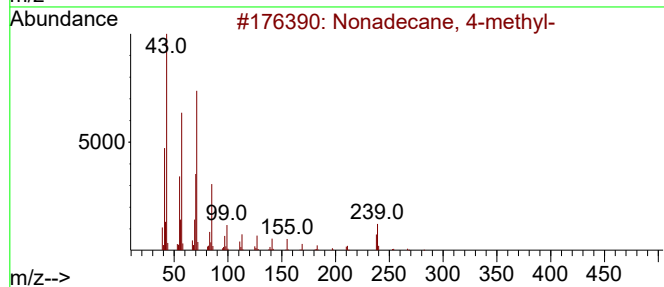
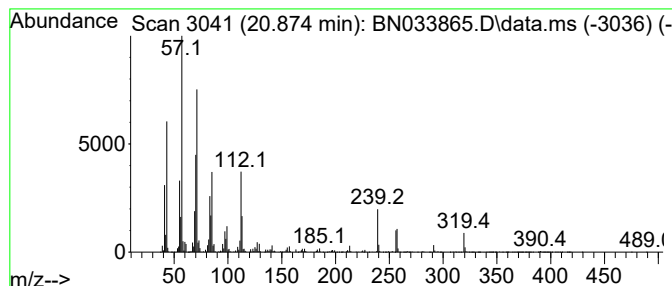
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 61 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.874	49.40 ng/ul	25930600	Chrysene-d12	21.139

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane, 4-methyl-	282	C20H42	025117-27-5	70
2			Carbonic acid, 2-ethylhexyl nony...	300	C18H36O3	1000383-13-6	62
3			Nonane, 4-methyl-5-propyl-	184	C13H28	062185-55-1	62
4			Oxalic acid, 2-ethylhexyl pentad...	412	C25H48O4	1000309-39-8	58
5			Heptadecane	240	C17H36	000629-78-7	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DDCC1

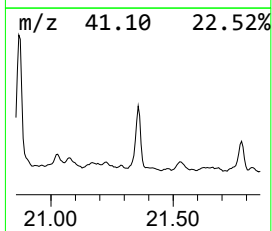
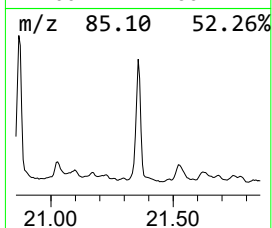
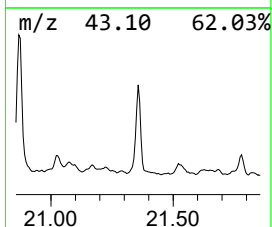
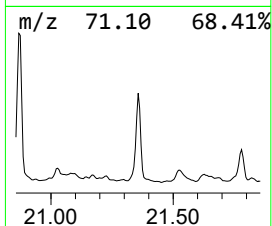
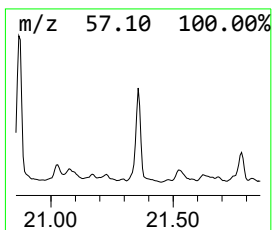
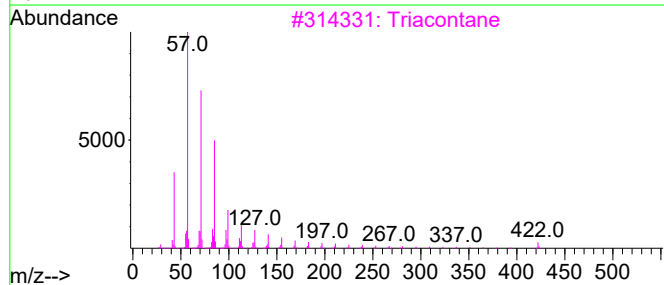
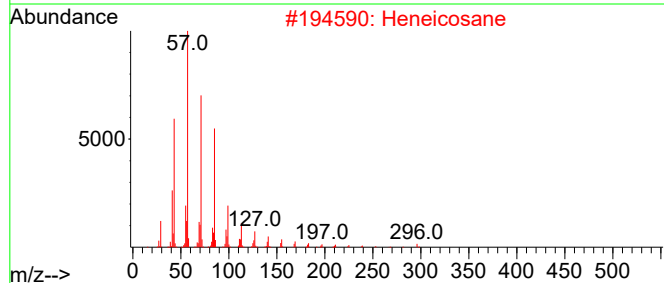
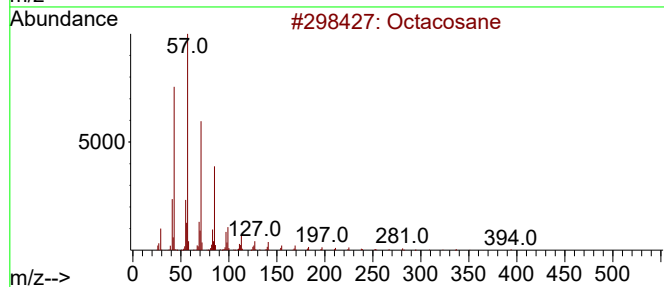
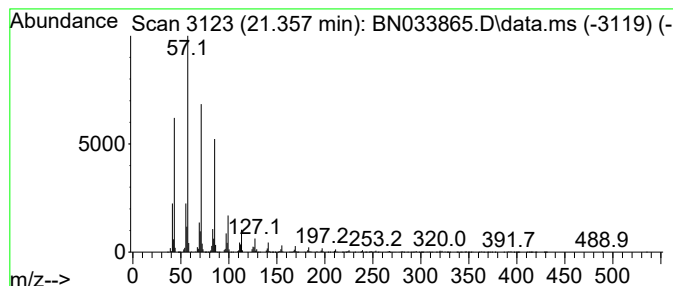
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 62 (DEL) Alkane: Straight-Chai... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.357	20.00 ng/ul	10499200	Chrysene-d12	21.139

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	99
2			Heneicosane	296	C21H44	000629-94-7	91
3			Triacontane	422	C30H62	000638-68-6	91
4			Tetracosane	338	C24H50	000646-31-1	91
5			Nonadecane	268	C19H40	000629-92-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DDCC1

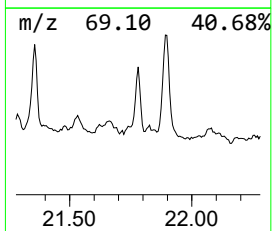
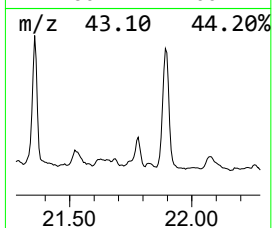
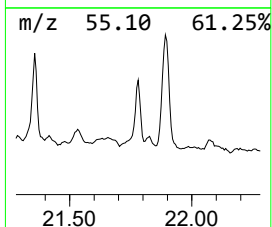
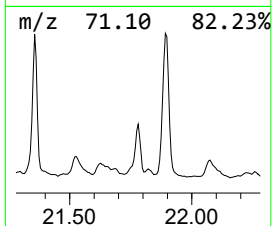
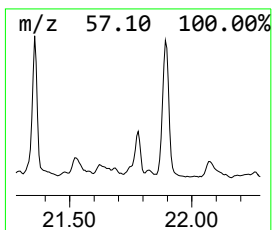
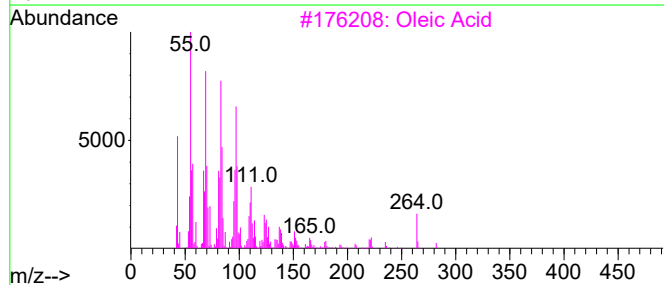
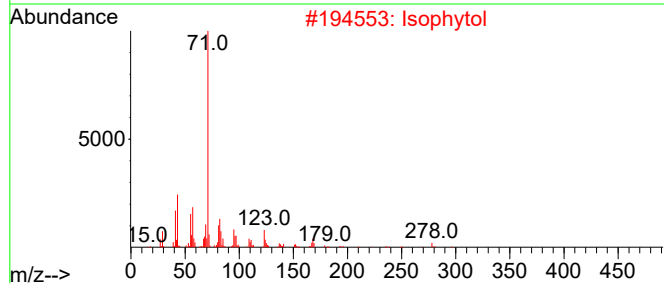
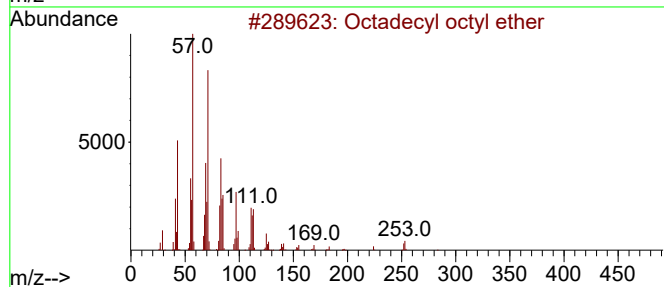
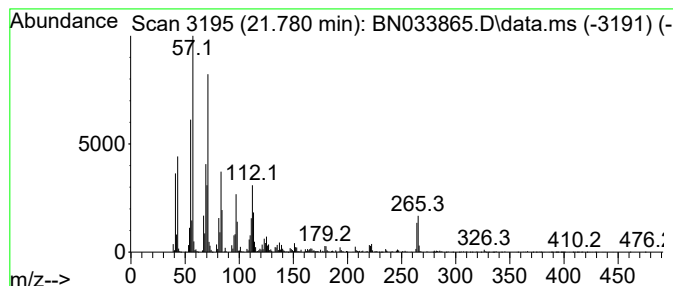
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 63 Octadecyl octyl ether Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.780	10.36 ng/ul	5440120	Chrysene-d12	21.139

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecyl octyl ether	382	C26H54O	1000406-38-7	64
2			Isophytol	296	C20H40O	000505-32-8	38
3			Oleic Acid	282	C18H34O2	000112-80-1	30
4			9-Octadecene, (E)-	252	C18H36	007206-25-9	25
5			5-Octadecene, (E)-	252	C18H36	007206-21-5	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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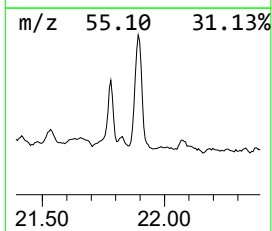
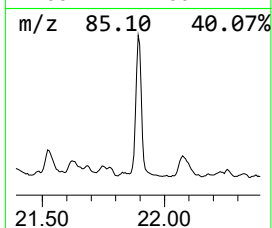
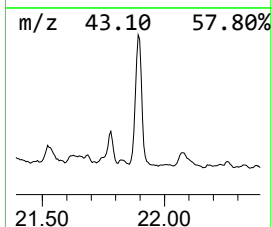
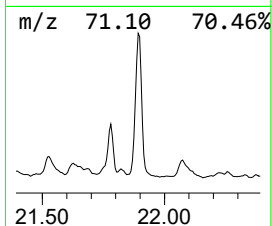
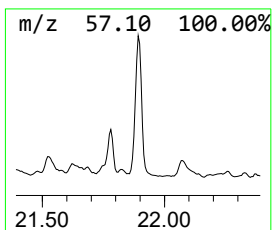
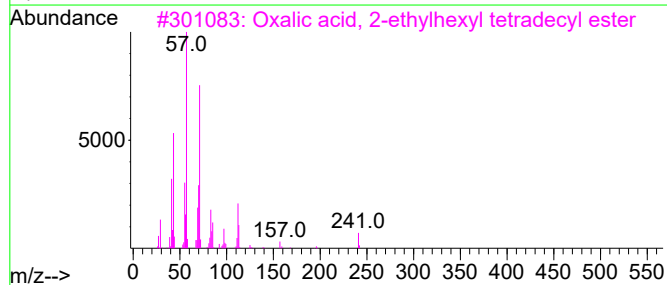
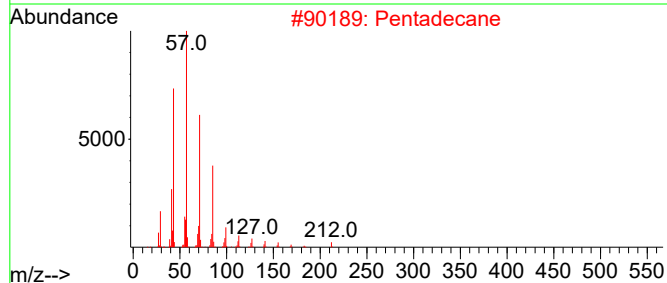
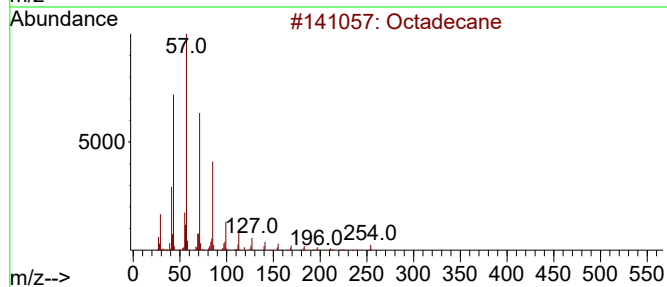
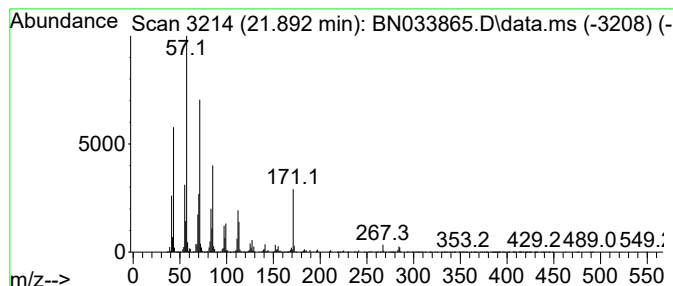
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 64 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.892	32.72 ng/ul	17177000	Chrysene-d12	21.139

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	90
2			Pentadecane	212	C15H32	000629-62-9	70
3			Oxalic acid, 2-ethylhexyl tetrad...	398	C24H46O4	1000309-39-7	62
4			Tetratetracontane	619	C44H90	007098-22-8	62
5			Decane, 2-methyl-	156	C11H24	006975-98-0	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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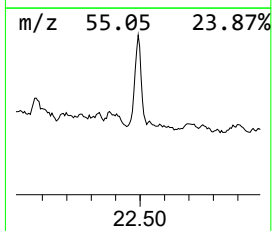
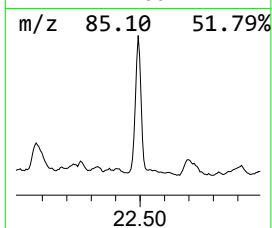
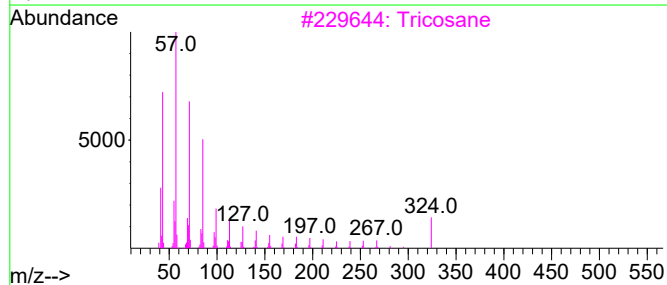
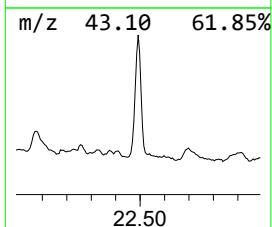
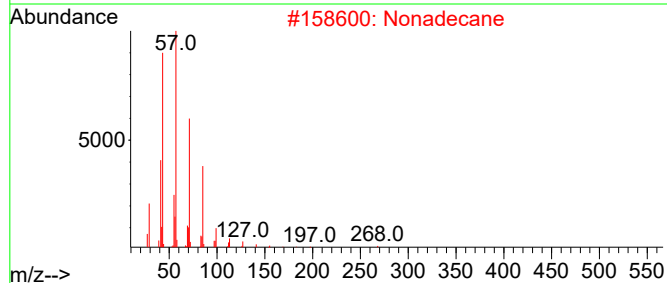
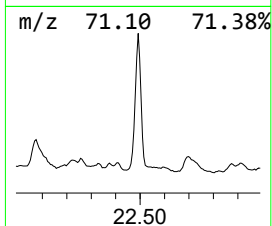
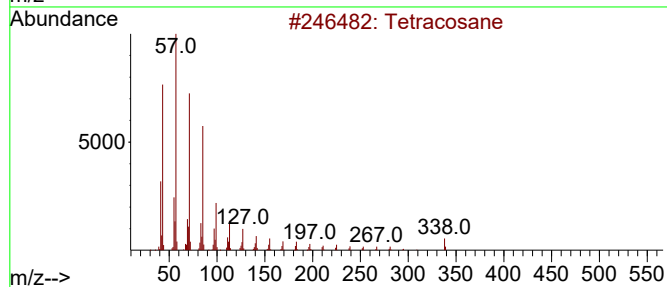
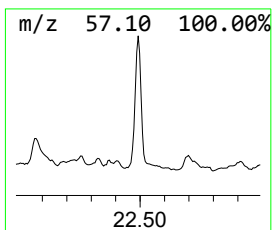
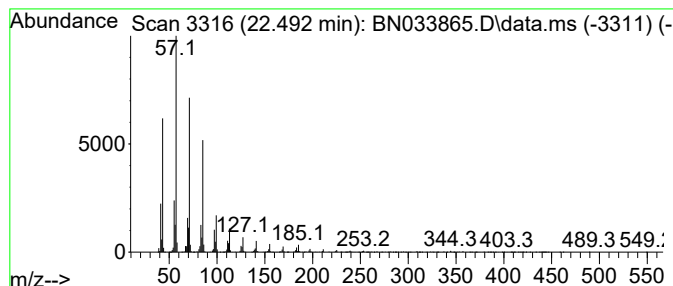
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 65 (DEL) Alkane: Straight-Chai... Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.492	13.56 ng/ul	7118750	Chrysene-d12	21.139

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	98
2			Nonadecane	268	C19H40	000629-92-5	95
3			Tricosane	324	C23H48	000638-67-5	95
4			2-Methylheptacosane	394	C28H58	001561-00-8	94
5			Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DDCC1

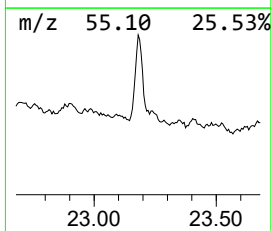
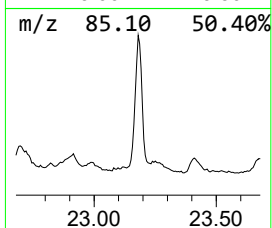
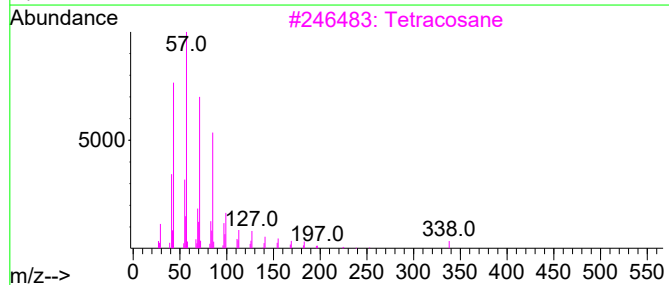
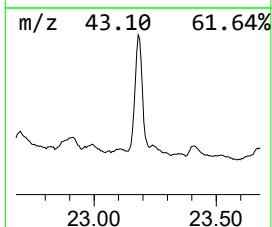
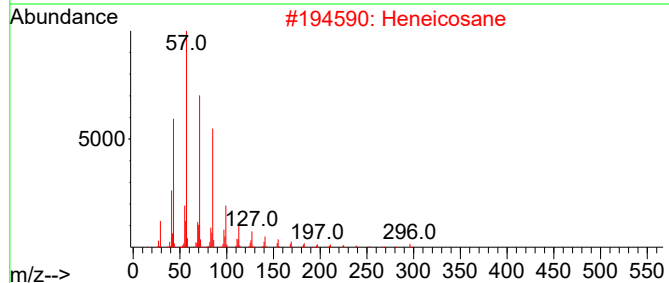
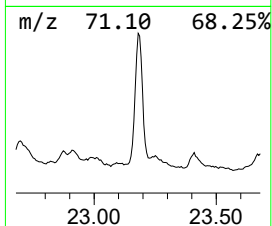
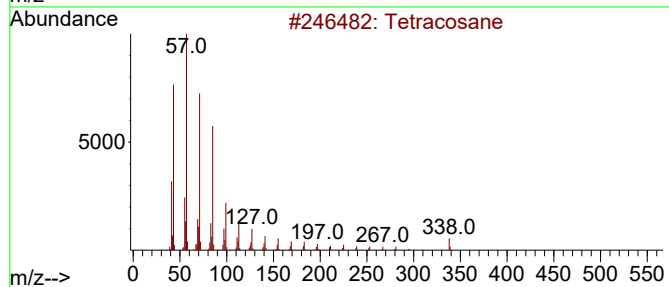
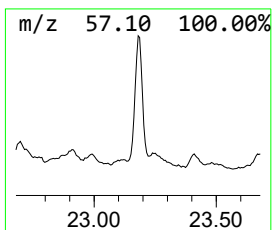
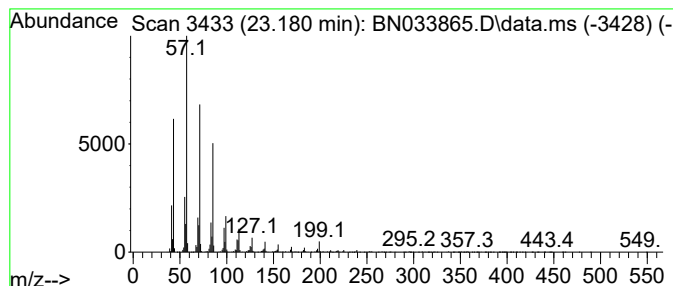
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 66 (DEL) Alkane: Straight-Chai... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.180	19.77 ng/ul	6077170	Perylene-d12	23.927

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	96
2			Heneicosane	296	C21H44	000629-94-7	91
3			Tetracosane	338	C24H50	000646-31-1	91
4			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5			Tetracosane	338	C24H50	000646-31-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

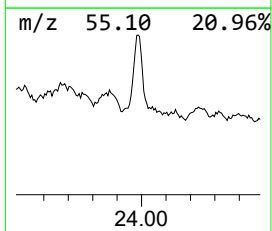
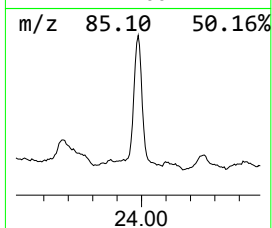
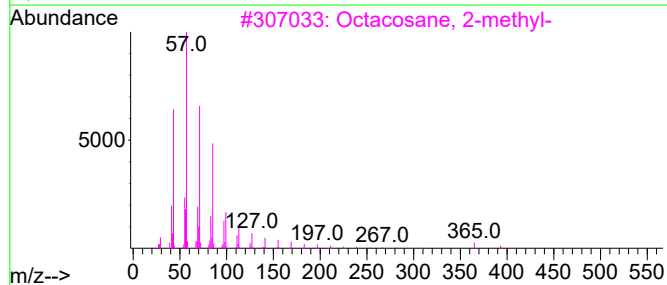
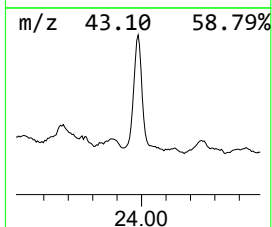
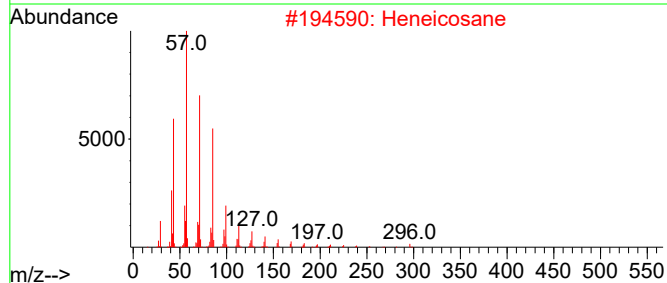
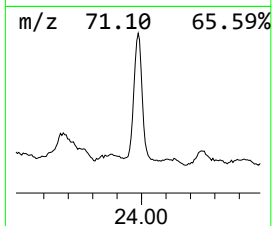
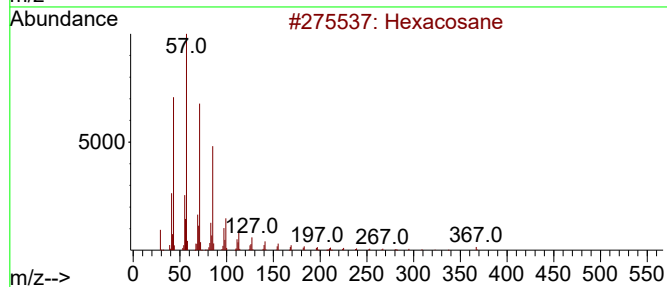
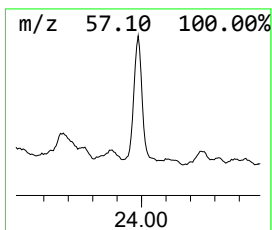
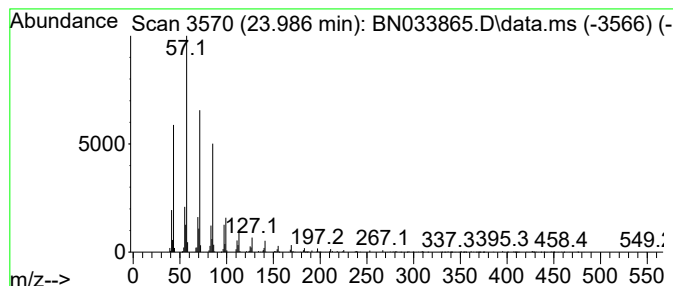
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Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 67 (DEL) Alkane: Straight-Chai... Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.986	14.41 ng/ul	4429160	Perylene-d12	23.927	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	91
2	Heneicosane	296	C21H44	000629-94-7	91
3	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
4	Octacosane	394	C28H58	000630-02-4	91
5	Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091124\
Data File : BN033865.D
Acq On : 11 Sep 2024 17:35
Operator : JU/RC
Sample : P3878-18
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	5.569	11.6	ng/ul	4487340	1	7.504	7768310	20.0
(DEL) Alkane: S...	6.581	12.2	ng/ul	4723880	1	7.504	7768310	20.0
2-Heptanone, 4,...	6.957	17.5	ng/ul	6807540	1	7.504	7768310	20.0
(DEL) Alkane: S...	7.157	72.7	ng/ul	28245400	1	7.504	7768310	20.0
1-Isobutoxy-2-e...	7.587	17.6	ng/ul	6817290	1	7.504	7768310	20.0
(DEL) Alkane: C...	7.775	16.7	ng/ul	6469150	1	7.504	7768310	20.0
Cyclohexanone, ...	7.934	22.6	ng/ul	8760370	1	7.504	7768310	20.0
unknown-01	7.993	26.7	ng/ul	10376700	1	7.504	7768310	20.0
Benzene, 1-meth...	8.046	25.2	ng/ul	9804530	1	7.504	7768310	20.0
(DEL) Alkane: S...	8.104	13.8	ng/ul	5353230	1	7.504	7768310	20.0
Benzene, 4-ethy...	8.140	14.6	ng/ul	5665160	1	7.504	7768310	20.0
(DEL) Alkane: S...	8.175	20.2	ng/ul	7843930	1	7.504	7768310	20.0
Benzene, 1-meth...	8.298	17.0	ng/ul	6595080	1	7.504	7768310	20.0
Benzene, 2-ethy...	8.451	10.3	ng/ul	4004300	1	7.504	7768310	20.0
o-Cymene	8.498	36.5	ng/ul	14165500	1	7.504	7768310	20.0
(DEL) Alkane: S...	8.740	57.9	ng/ul	22506500	1	7.504	7768310	20.0
unknown-02	8.845	18.8	ng/ul	7312680	1	7.504	7768310	20.0
2H-Pyran-2-carb...	12.392	8.5	ng/ul	8983920	3	14.145	21052100	20.0
(DEL) Alkane: C...	12.486	5.5	ng/ul	5775650	3	14.145	21052100	20.0
Bicyclo[4.1.0]h...	12.728	14.0	ng/ul	14783500	3	14.145	21052100	20.0
2,6-Pyridinedia...	13.210	14.6	ng/ul	15313400	3	14.145	21052100	20.0
Naphthalene, 1,...	13.304	6.2	ng/ul	6534760	3	14.145	21052100	20.0
Naphthalene, 2,...	13.469	8.0	ng/ul	8363060	3	14.145	21052100	20.0
Naphthalene, 1,...	13.516	10.2	ng/ul	10685900	3	14.145	21052100	20.0
(DEL) Alkane: S...	13.657	4.3	ng/ul	4482250	3	14.145	21052100	20.0
1-(benzo[d][1,3...	13.904	7.4	ng/ul	7816270	3	14.145	21052100	20.0
Sulfurous acid,...	14.086	20.0	ng/ul	21052100	3	14.145	21052100	20.0
unknown-03	14.251	6.0	ng/ul	6363900	3	14.145	21052100	20.0
9-Methyl-5-octa...	14.310	15.7	ng/ul	16566100	3	14.145	21052100	20.0
unknown-04	14.704	9.0	ng/ul	9501650	3	14.145	21052100	20.0
3,4-Dimethyl(1H...	14.933	21.7	ng/ul	22868400	3	14.145	21052100	20.0
(DEL) Alkane: S...	15.045	20.0	ng/ul	21025100	3	14.145	21052100	20.0
(DEL) Alkane: S...	15.445	16.9	ng/ul	17757400	3	14.145	21052100	20.0
(DEL) Alkane: S...	15.598	8.5	ng/ul	4707280	4	16.939	11121100	20.0
unknown-05	15.804	16.8	ng/ul	9327280	4	16.939	11121100	20.0
(DEL) Alkane: S...	15.910	36.5	ng/ul	20305600	4	16.939	11121100	20.0
(DEL) Alkane: S...	16.757	28.3	ng/ul	15744800	4	16.939	11121100	20.0
unknown-06	16.833	7.2	ng/ul	4020850	4	16.939	11121100	20.0
unknown-07	17.127	9.3	ng/ul	5173180	4	16.939	11121100	20.0
unknown-08	17.245	24.1	ng/ul	13387700	4	16.939	11121100	20.0
(DEL) Alkane: S...	17.445	31.8	ng/ul	17668400	4	16.939	11121100	20.0
2-Methyldibenzo...	17.498	9.9	ng/ul	5523820	4	16.939	11121100	20.0
n-Hexadecanoic ...	17.863	36.1	ng/ul	20102000	4	16.939	11121100	20.0
(DEL) Alkane: S...	18.133	27.0	ng/ul	14994300	4	16.939	11121100	20.0
(DEL) Alkane: S...	18.774	22.6	ng/ul	12580700	4	16.939	11121100	20.0
Octadecanoic acid	19.139	14.3	ng/ul	7505790	5	21.139	10499200	20.0
(DEL) Alkane: S...	19.898	23.6	ng/ul	12377800	5	21.139	10499200	20.0
(DEL) Alkane: S...	20.398	18.7	ng/ul	9822840	5	21.139	10499200	20.0
(DEL) Alkane: S...	20.874	49.4	ng/ul	25930600	5	21.139	10499200	20.0
(DEL) Alkane: S...	21.357	20.0	ng/ul	10499200	5	21.139	10499200	20.0
Octadecyl octyl...	21.780	10.4	ng/ul	5440120	5	21.139	10499200	20.0
(DEL) Alkane: S...	21.892	32.7	ng/ul	17177000	5	21.139	10499200	20.0

(DEL) Alkane: S... 22.492 13.6 ng/ul 7118750 5 21.139 10499200 20.0
(DEL) Alkane: S... 23.180 19.8 ng/ul 6077170 6 23.927 6149170 20.0
(DEL) Alkane: S... 23.986 14.4 ng/ul 4429160 6 23.927 6149170 20.0

Instrument :
BNA_N
ClientSampleId :
DDCC1