

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100224\
 Data File : BN034363.D
 Acq On : 03 Oct 2024 09:58
 Operator : RC/JU
 Sample : P4018-04DL 20X
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DDCA2DL

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: BN034363.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.693	132	137	142	rVB	37934	53065	0.89%	0.170%
2	5.446	429	435	444	rVV2	121234	186802	3.15%	0.597%
3	5.975	521	525	530	rBV	36310	46875	0.79%	0.150%
4	6.046	530	537	541	rVV4	27125	61999	1.04%	0.198%
5	6.311	575	582	587	rBV4	21368	46701	0.79%	0.149%
6	6.405	594	598	602	rVB2	49919	73536	1.24%	0.235%
7	6.458	602	607	610	rBV	64913	91574	1.54%	0.293%
8	6.493	610	613	617	rVV	59657	85599	1.44%	0.274%
9	6.564	617	625	631	rVV4	81559	199303	3.36%	0.637%
10	6.622	631	635	641	rVV	57214	87808	1.48%	0.281%
11	6.740	648	655	658	rVV2	24273	45267	0.76%	0.145%
12	6.781	658	662	666	rVV2	51389	81031	1.36%	0.259%
13	6.834	666	671	676	rVV3	31051	67193	1.13%	0.215%
14	6.922	682	686	698	rVV2	102320	188297	3.17%	0.602%
15	7.028	698	704	714	rVV2	368489	630918	10.62%	2.016%
16	7.381	757	764	771	rVV	810686	1259210	21.20%	4.025%
17	7.464	772	778	791	rVB2	260222	465225	7.83%	1.487%
18	7.617	800	804	809	rBV2	66626	107770	1.81%	0.344%
19	7.922	852	856	861	rVB2	75120	122315	2.06%	0.391%
20	7.975	861	865	869	rBV	44017	63045	1.06%	0.201%
21	8.017	869	872	875	rBV	47180	71418	1.20%	0.228%
22	8.046	875	877	883	rVB	60915	78326	1.32%	0.250%
23	8.152	890	895	897	rBV	59751	81159	1.37%	0.259%
24	8.328	920	925	929	rBV	33480	51557	0.87%	0.165%
25	8.375	929	933	940	rVV	35977	58822	0.99%	0.188%
26	8.475	940	950	954	rVV2	45034	93740	1.58%	0.300%
27	8.522	954	958	963	rVV2	23164	50621	0.85%	0.162%
28	8.611	968	973	980	rVB	262547	375877	6.33%	1.201%
29	8.681	980	985	988	rBV2	53766	92353	1.55%	0.295%
30	8.722	989	992	999	rVB5	33298	60146	1.01%	0.192%
31	8.805	999	1006	1011	rBV3	48366	92439	1.56%	0.295%
32	9.240	1073	1080	1086	rBV7	28013	72200	1.22%	0.231%
33	9.322	1086	1094	1098	rVV4	26817	56941	0.96%	0.182%
34	9.452	1106	1116	1120	rVB4	19758	51809	0.87%	0.166%
35	9.775	1167	1171	1179	rVB2	30928	56459	0.95%	0.180%
36	10.040	1209	1216	1224	rBV2	45783	102789	1.73%	0.329%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0.1 % of largest Peak

Max Peaks: 100

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	10.140	1224	1233	1240	rBV	1091574	1899613	31.98%	6.071%
38	10.281	1251	1257	1262	rBV4	77504	142467	2.40%	0.455%
39	10.534	1294	1300	1308	rBV	55303	93979	1.58%	0.300%
40	10.840	1346	1352	1357	rBV3	26923	47921	0.81%	0.153%

41	11.187	1406	1411	1417	rBV6	32711	81975	1.38%	0.262%
42	11.252	1417	1422	1428	rVB	119357	182640	3.08%	0.584%
43	11.428	1448	1452	1458	rVB2	49284	74894	1.26%	0.239%
44	11.599	1472	1481	1484	rBV	68297	132791	2.24%	0.424%
45	11.628	1484	1486	1491	rVB	50111	62768	1.06%	0.201%

46	11.681	1491	1495	1502	rBV4	28459	51507	0.87%	0.165%
47	11.857	1521	1525	1534	rVB	66601	106547	1.79%	0.341%
48	12.034	1546	1555	1559	rBV2	25580	61760	1.04%	0.197%
49	12.263	1588	1594	1598	rBV2	46959	77482	1.30%	0.248%
50	12.581	1643	1648	1656	rVB3	23991	43764	0.74%	0.140%

51	12.787	1674	1683	1689	rBV3	17795	43895	0.74%	0.140%
52	12.875	1689	1698	1705	rVV	88240	141347	2.38%	0.452%
53	13.193	1746	1752	1758	rBV4	26555	69930	1.18%	0.224%
54	13.352	1775	1779	1784	rVV2	39989	67538	1.14%	0.216%
55	13.404	1784	1788	1793	rVV3	43754	77201	1.30%	0.247%

56	13.457	1793	1797	1803	rVV	56430	86562	1.46%	0.277%
57	13.552	1805	1813	1816	rVV2	26274	50368	0.85%	0.161%
58	13.687	1829	1836	1839	rBV3	61650	103959	1.75%	0.332%
59	13.722	1839	1842	1846	rVV	73751	101456	1.71%	0.324%
60	13.969	1879	1884	1889	rBV	754071	986147	16.60%	3.152%

61	14.034	1889	1895	1903	rVV	1460840	2100965	35.37%	6.715%
62	14.193	1918	1922	1927	rVB	177715	235602	3.97%	0.753%
63	14.269	1927	1935	1940	rBV5	31058	78912	1.33%	0.252%
64	14.375	1948	1953	1957	rVV5	21943	45818	0.77%	0.146%
65	14.416	1957	1960	1968	rVB9	24119	60529	1.02%	0.193%

66	14.675	1999	2004	2009	rBV	315726	453435	7.63%	1.449%
67	14.763	2015	2019	2023	rVB	57606	70229	1.18%	0.224%
68	14.816	2023	2028	2035	rBV	190918	282548	4.76%	0.903%
69	14.934	2040	2048	2052	rVB2	108551	182374	3.07%	0.583%
70	15.040	2061	2066	2070	rBV	92604	130886	2.20%	0.418%

71	15.075	2070	2072	2080	rVB7	31774	46336	0.78%	0.148%
72	15.157	2080	2086	2094	rBV	52153	99251	1.67%	0.317%
73	15.340	2110	2117	2122	rBV2	31772	68831	1.16%	0.220%
74	15.422	2127	2131	2138	rVB4	35695	59374	1.00%	0.190%
75	15.698	2171	2178	2184	rBV3	75438	108237	1.82%	0.346%

76	15.798	2190	2195	2198	rBV	115170	156954	2.64%	0.502%
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 Title : SVOA CALIBRATION

77	15.875	2203	2208	2213	rVV	79701	113746	1.92%	0.364%
78	16.128	2247	2251	2261	rVB6	21624	45496	0.77%	0.145%
79	16.451	2299	2306	2319	rBV	4291132	5939297	100.00%	18.983%
80	16.593	2326	2330	2336	rBV2	82470	118533	2.00%	0.379%
81	16.645	2336	2339	2345	rVB2	34819	52452	0.88%	0.168%
82	16.793	2357	2364	2376	rBV	1773473	2494243	42.00%	7.972%
83	16.893	2376	2381	2388	rVV	98161	175262	2.95%	0.560%
84	17.334	2450	2456	2460	rVV	70304	96014	1.62%	0.307%
85	17.381	2460	2464	2469	rVV3	21380	44099	0.74%	0.141%
86	17.504	2480	2485	2493	rBV2	54408	86775	1.46%	0.277%
87	18.028	2568	2574	2578	rBV	50192	72466	1.22%	0.232%
88	18.669	2679	2683	2685	rVV3	45885	64831	1.09%	0.207%
89	18.692	2685	2687	2693	rVV5	47203	52670	0.89%	0.168%
90	19.163	2763	2767	2769	rBV2	31645	42780	0.72%	0.137%
91	19.204	2770	2774	2780	rVV	108475	158909	2.68%	0.508%
92	19.798	2872	2875	2881	rVB2	35983	43168	0.73%	0.138%
93	19.869	2883	2887	2891	rBV	54717	69079	1.16%	0.221%
94	20.504	2991	2995	3000	rBV	105963	130615	2.20%	0.417%
95	20.769	3037	3040	3044	rBV	54446	61640	1.04%	0.197%
96	21.028	3079	3084	3092	rBV	2044103	2825538	47.57%	9.031%
97	21.104	3093	3097	3103	rVB	160102	217768	3.67%	0.696%
98	21.751	3203	3207	3218	rVB2	165655	316008	5.32%	1.010%
99	22.480	3327	3331	3341	rVB2	147204	271545	4.57%	0.868%
100	23.739	3536	3545	3553	rBV	1489246	3415966	57.51%	10.918%

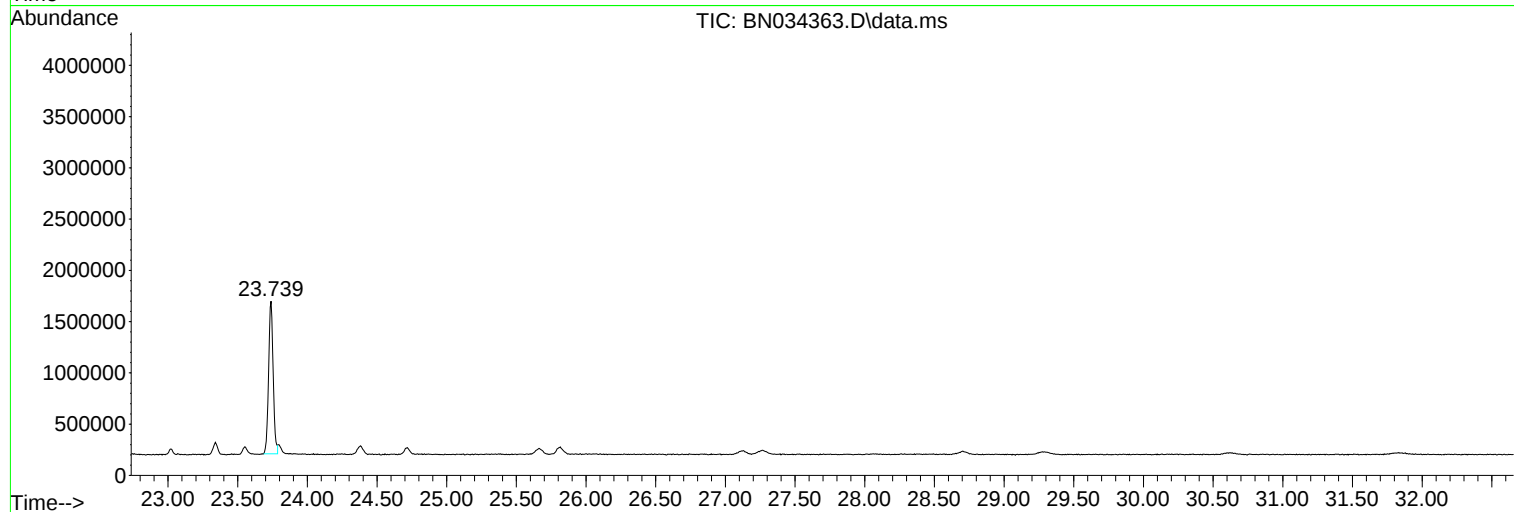
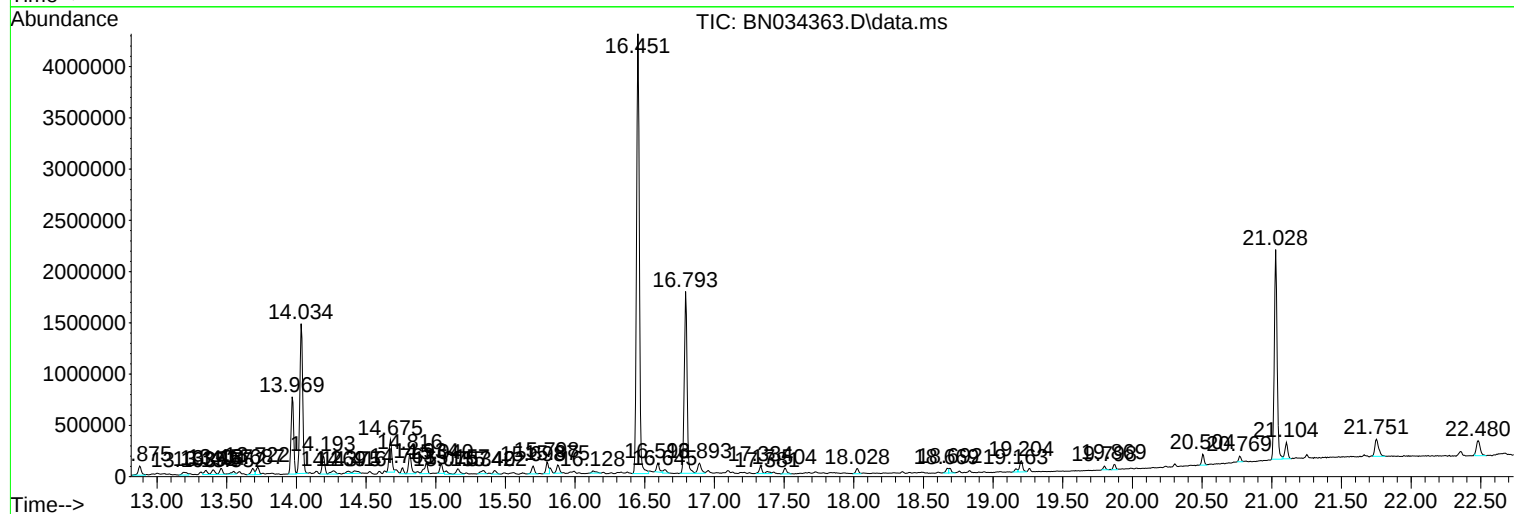
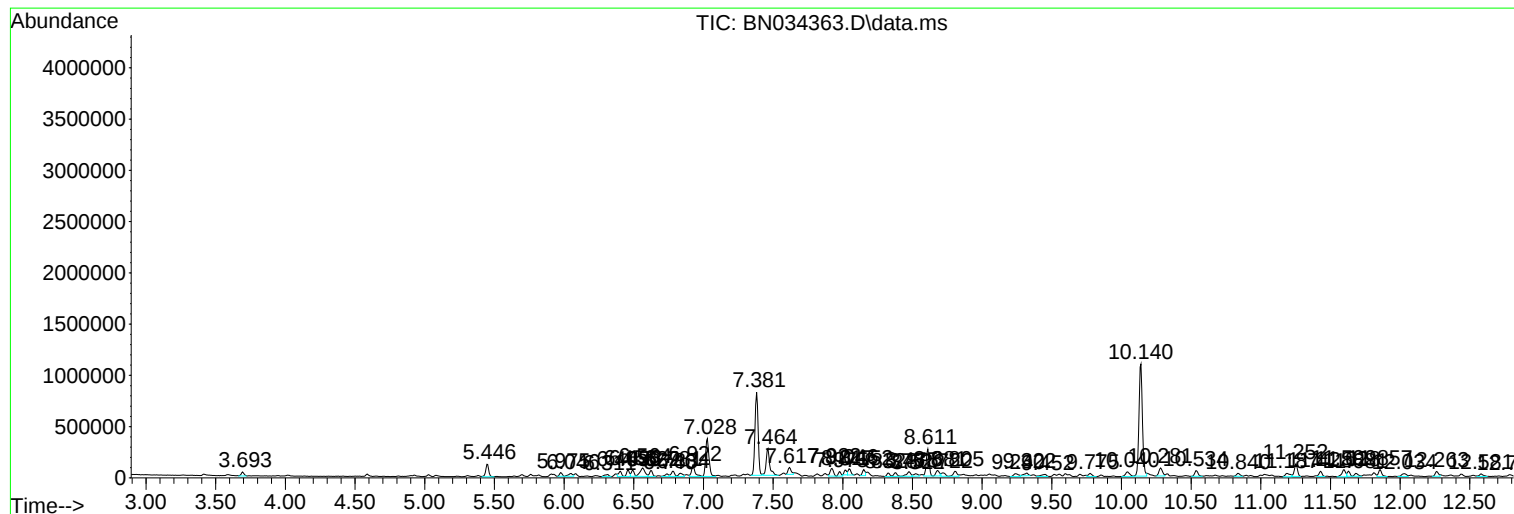
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ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DDCA2DL

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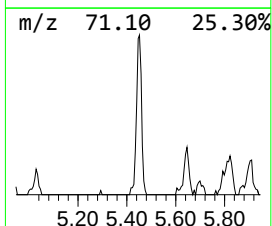
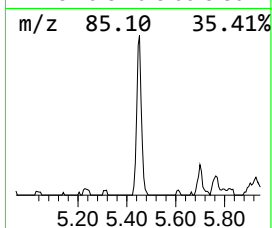
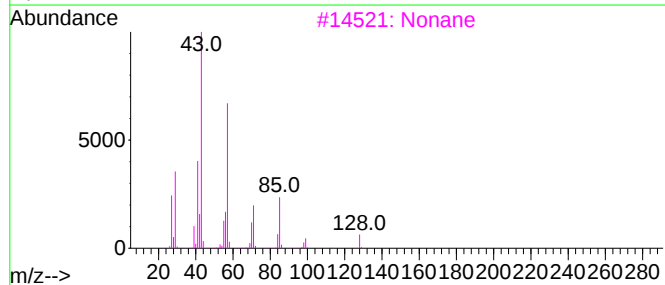
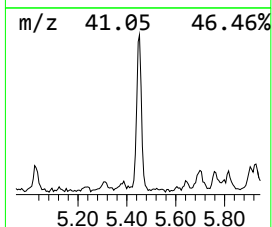
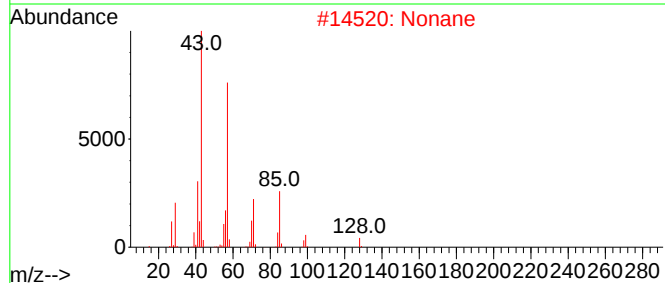
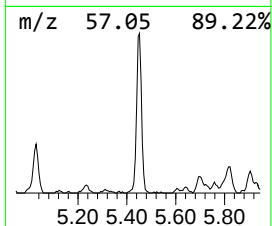
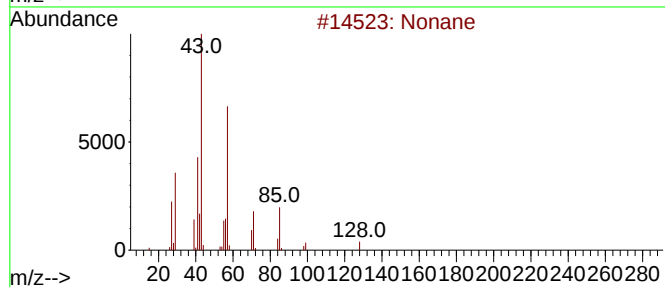
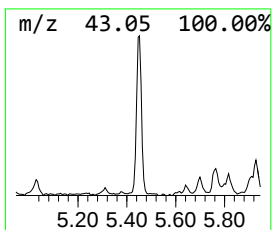
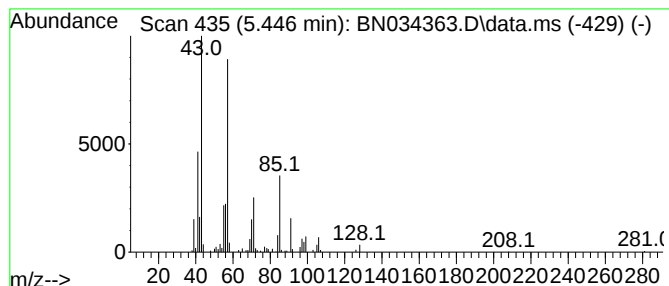
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.446	2.97 ng/ul	186802	1,4-Dichlorobenzene-d4	7.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane			128	C9H20	000111-84-2	90
2	Nonane			128	C9H20	000111-84-2	81
3	Nonane			128	C9H20	000111-84-2	59
4	Octane			114	C8H18	000111-65-9	47
5	Decane			142	C10H22	000124-18-5	47



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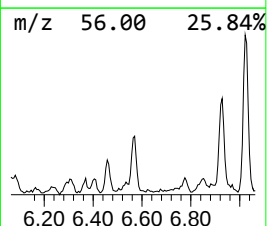
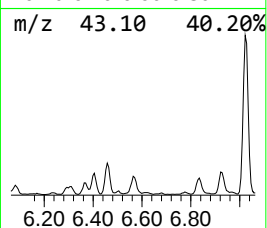
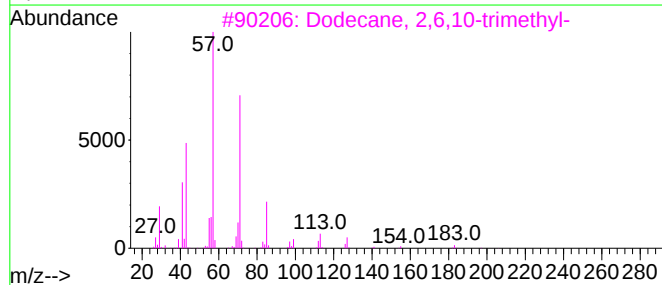
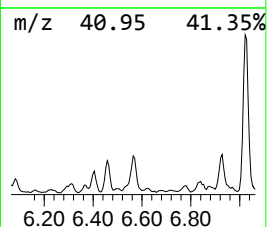
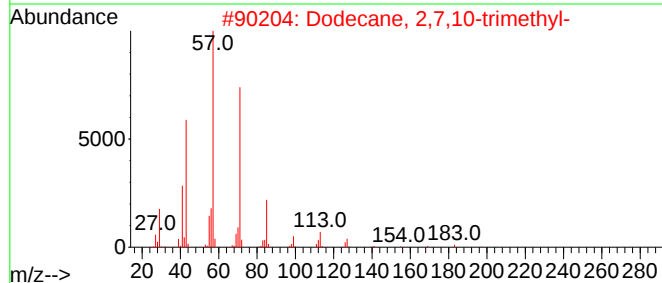
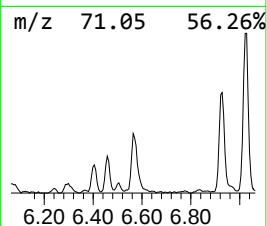
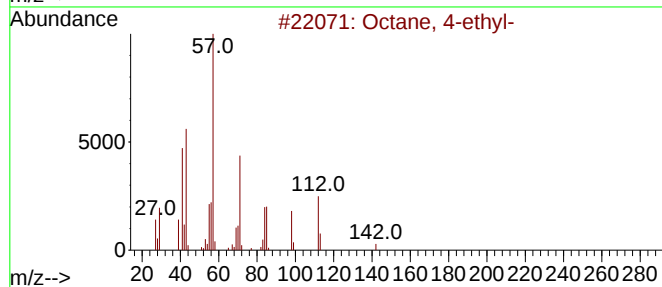
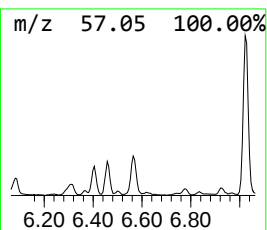
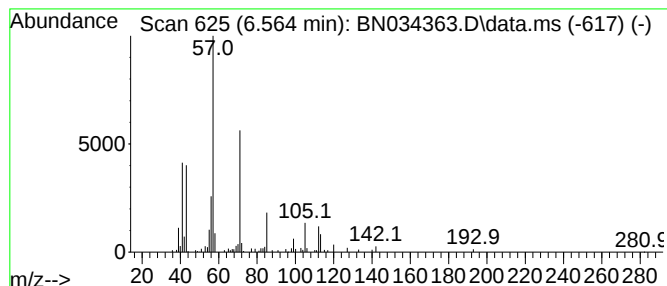
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.564	3.17 ng/ul	199303	1,4-Dichlorobenzene-d4	7.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 4-ethyl-			142	C10H22	015869-86-0	64
2	Dodecane, 2,7,10-trimethyl-			212	C15H32	074645-98-0	64
3	Dodecane, 2,6,10-trimethyl-			212	C15H32	003891-98-3	59
4	Octane, 3,5-dimethyl-			142	C10H22	015869-93-9	59
5	Octane, 2,6-dimethyl-			142	C10H22	002051-30-1	59



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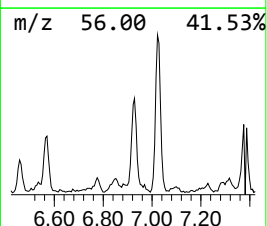
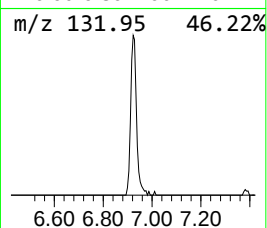
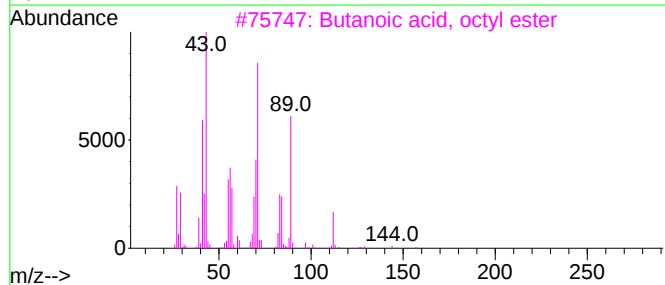
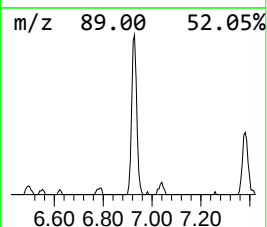
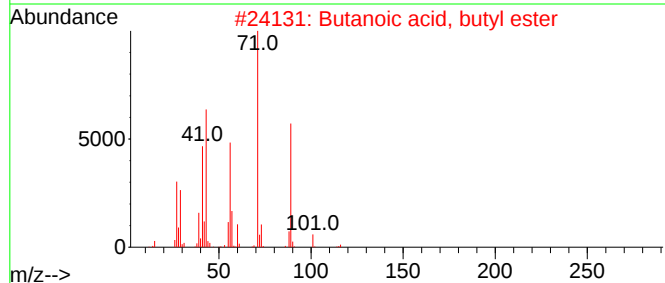
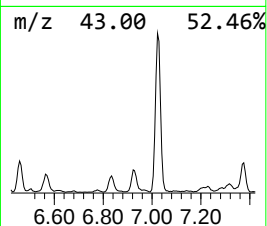
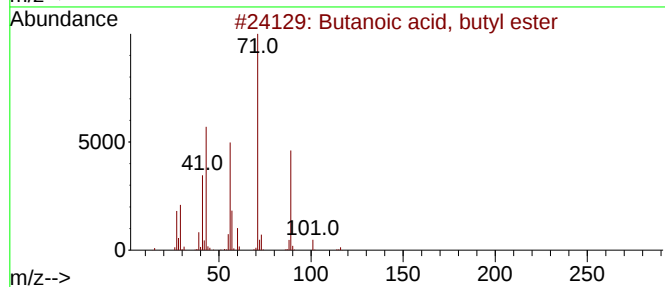
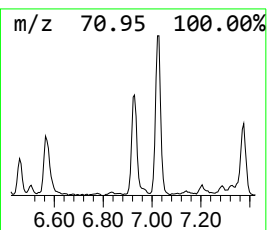
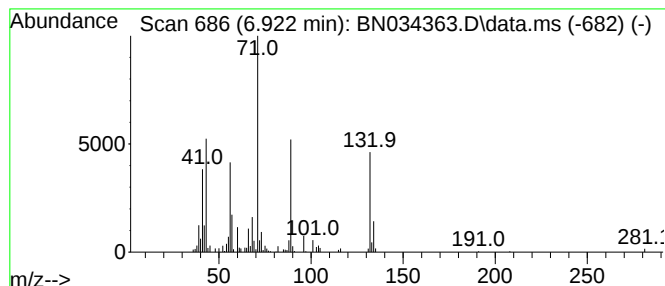
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Butanoic acid, butyl ester Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.922	2.99 ng/ul	188297	1,4-Dichlorobenzene-d4	7.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	58
2			Butanoic acid, butyl ester	144	C8H16O2	000109-21-7	58
3			Butanoic acid, octyl ester	200	C12H24O2	000110-39-4	53
4			Butanoic acid, 2-methylpropyl ester	144	C8H16O2	000539-90-2	47
5			Butanoic acid, 2-methylpropyl ester	144	C8H16O2	000539-90-2	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100224\
Data File : BN034363.D
Acq On : 03 Oct 2024 09:58
Operator : RC/JU
Sample : P4018-04DL 20X
Misc :
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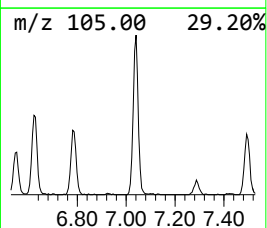
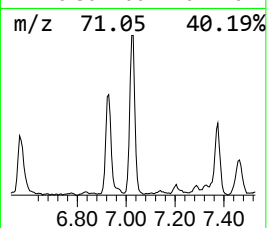
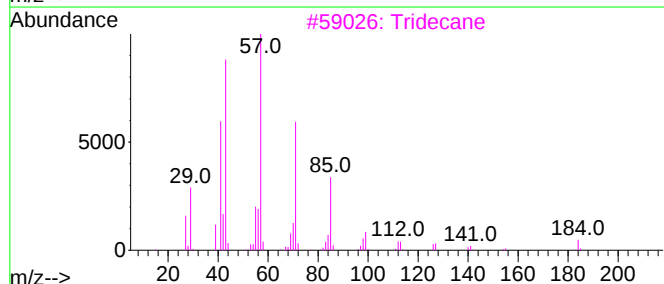
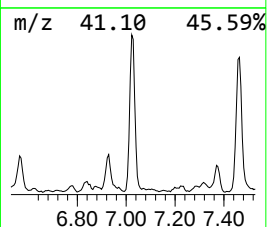
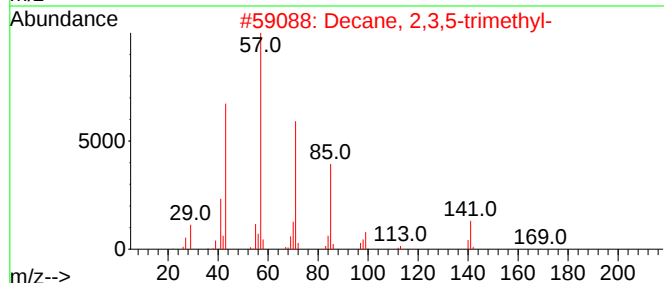
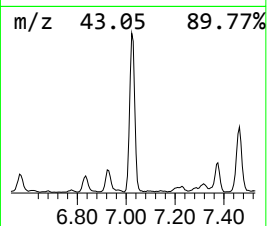
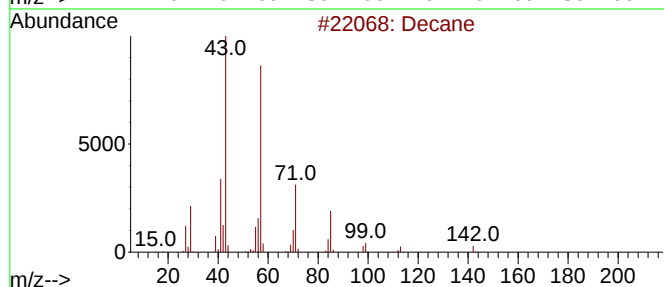
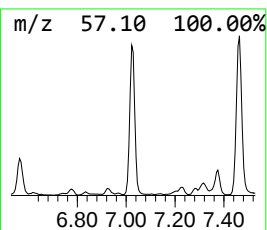
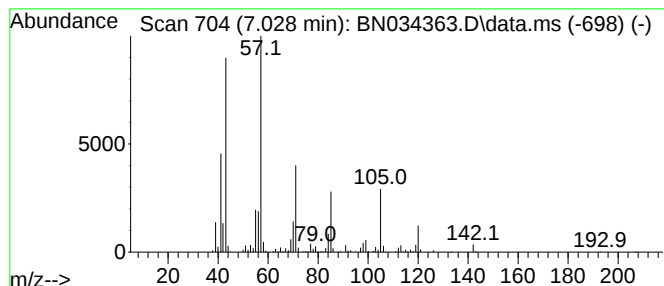
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Quant Title : SVOA CALIBRATION

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.
7.028	10.02 ng/ul	630918	1,4-Dichlorobenzene-d4			7.381
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Decane		142	C10H22	000124-18-5	89
2	Decane, 2,3,5-trimethyl-		184	C13H28	062238-11-3	58
3	Tridecane		184	C13H28	000629-50-5	58
4	Dodecane		170	C12H26	000112-40-3	53
5	Tridecane		184	C13H28	000629-50-5	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100224\
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Acq On : 03 Oct 2024 09:58
Operator : RC/JU
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ClientSampleId :
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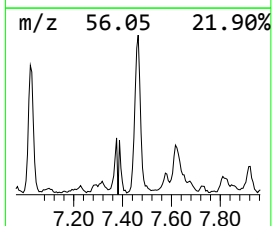
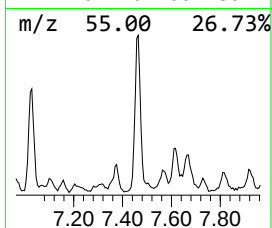
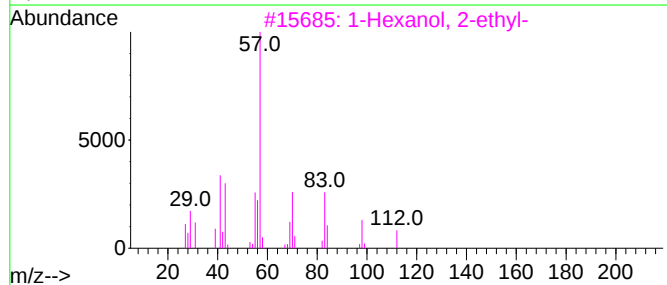
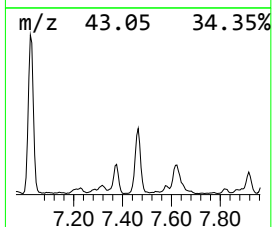
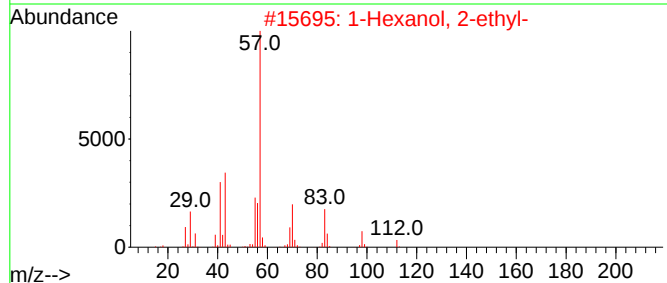
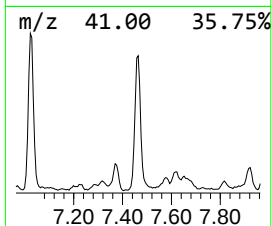
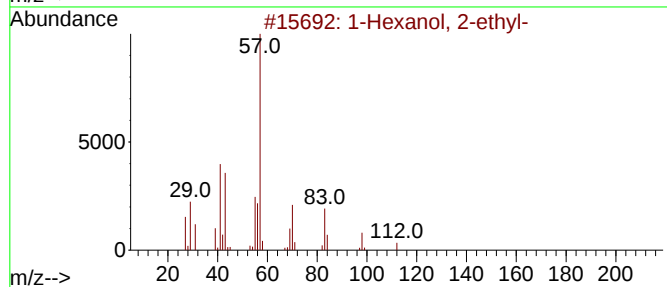
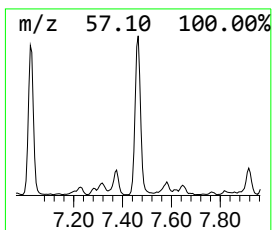
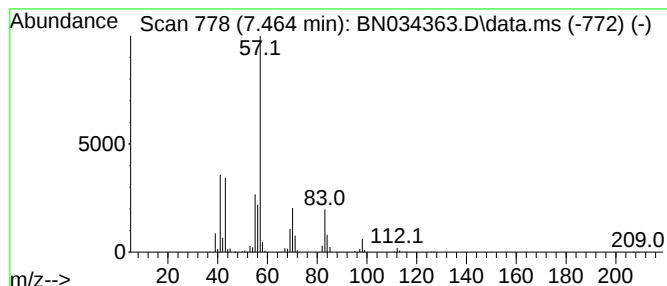
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1-Hexanol, 2-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.464	7.39 ng/ul	465225	1,4-Dichlorobenzene-d4	7.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-		130	C8H18O	000104-76-7	83	
2	1-Hexanol, 2-ethyl-		130	C8H18O	000104-76-7	83	
3	1-Hexanol, 2-ethyl-		130	C8H18O	000104-76-7	83	
4	1-Hexanol, 2-ethyl-		130	C8H18O	000104-76-7	74	
5	1-Hexanol, 2-ethyl-		130	C8H18O	000104-76-7	64	



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100224\
Data File : BN034363.D
Acq On : 03 Oct 2024 09:58
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Sample : P4018-04DL 20X
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
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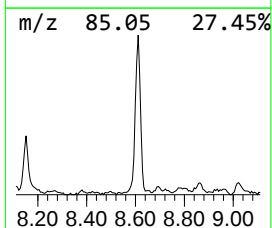
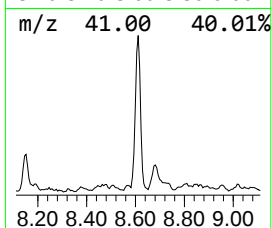
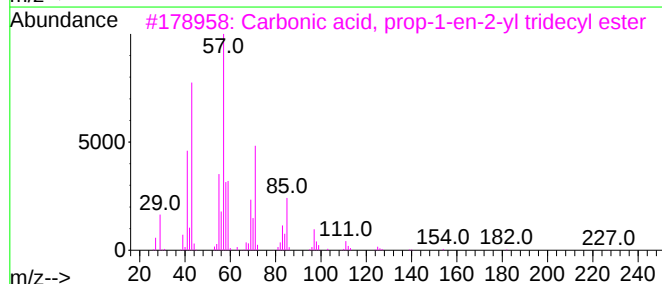
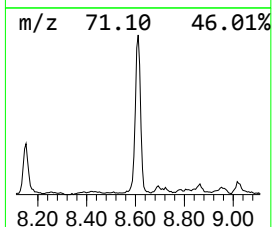
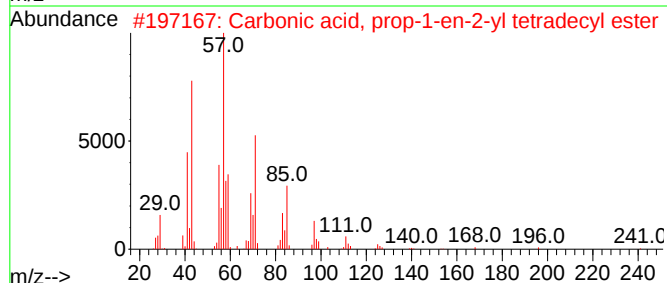
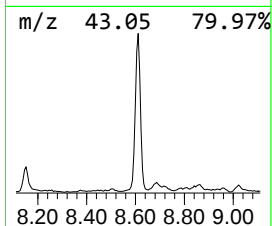
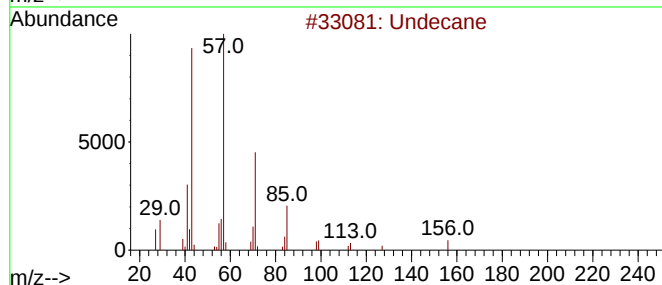
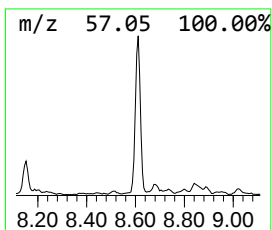
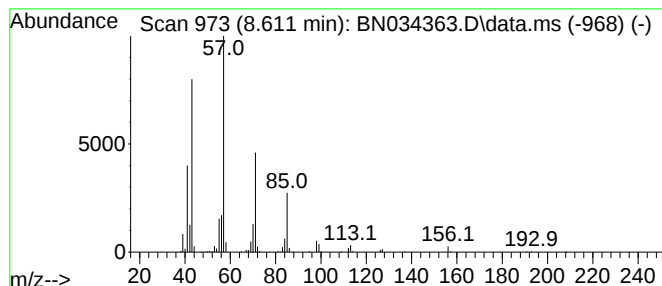
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.611	5.97 ng/ul	375877	1,4-Dichlorobenzene-d4	7.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane	156	C11H24	001120-21-4	91
2			Carbonic acid, prop-1-en-2-yl te...	298	C18H34O3	1000382-90-9	90
3			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	83
4			Tridecane	184	C13H28	000629-50-5	78
5			Undecane	156	C11H24	001120-21-4	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100224\
Data File : BN034363.D
Acq On : 03 Oct 2024 09:58
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Sample : P4018-04DL 20X
Misc :
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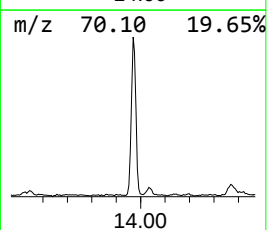
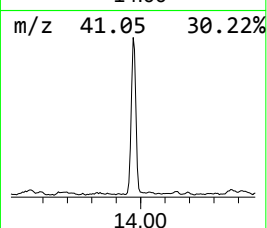
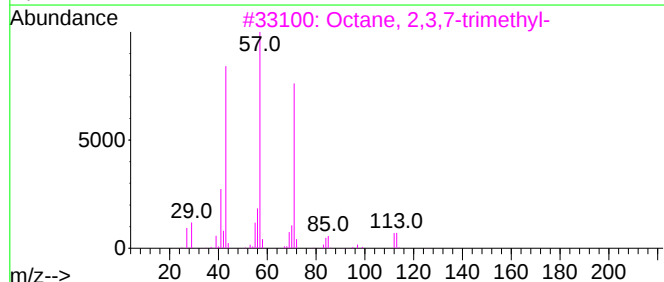
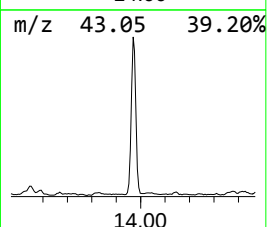
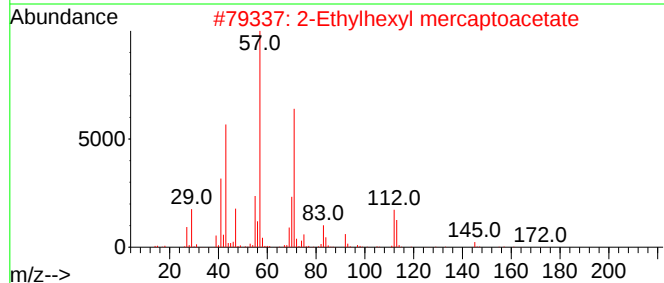
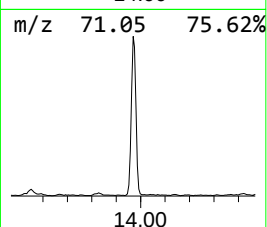
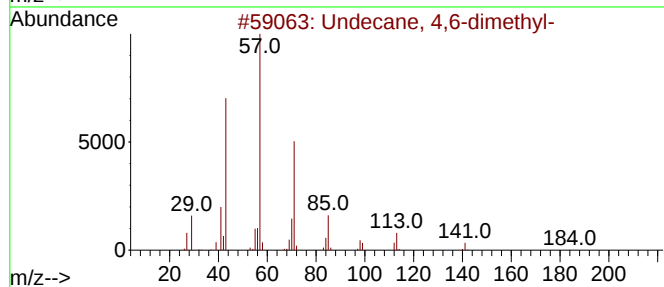
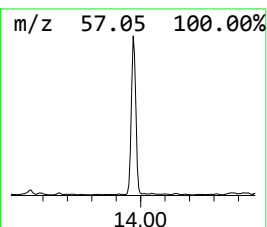
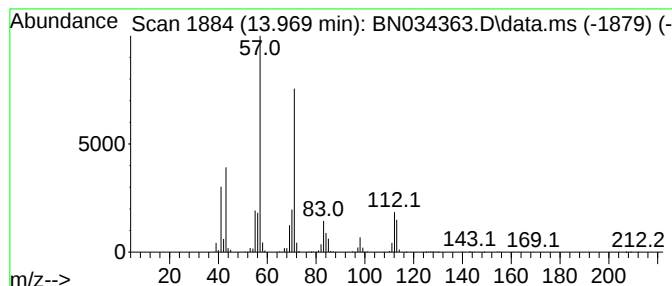
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.969	9.39 ng/ul	986147	Acenaphthene-d10	14.034

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 4,6-dimethyl-	184	C13H28	017312-82-2	78
2			2-Ethylhexyl mercaptoacetate	204	C10H20O2S	007659-86-1	78
3			Octane, 2,3,7-trimethyl-	156	C11H24	062016-34-6	72
4			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	64
5			Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1010309-19-2	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100224\
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Acq On : 03 Oct 2024 09:58
Operator : RC/JU
Sample : P4018-04DL 20X
Misc :
ALS Vial : 50 Sample Multiplier: 1

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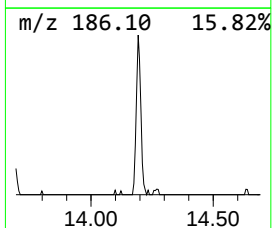
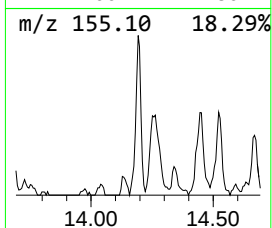
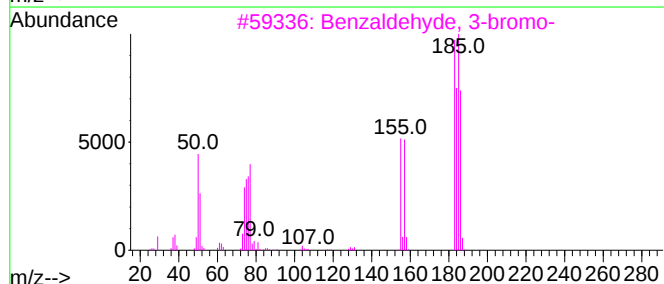
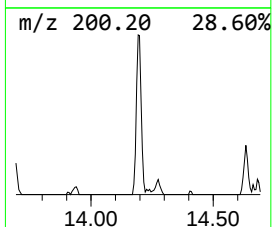
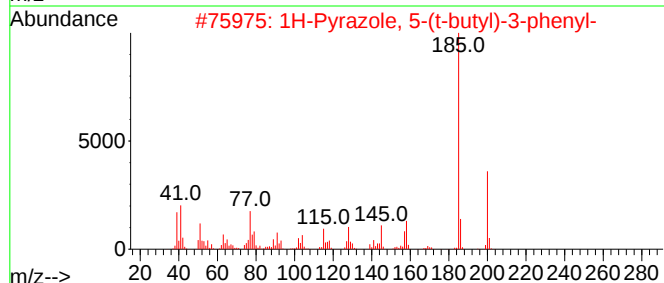
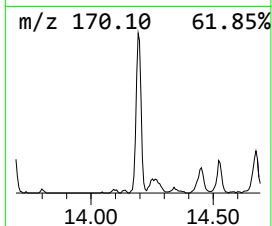
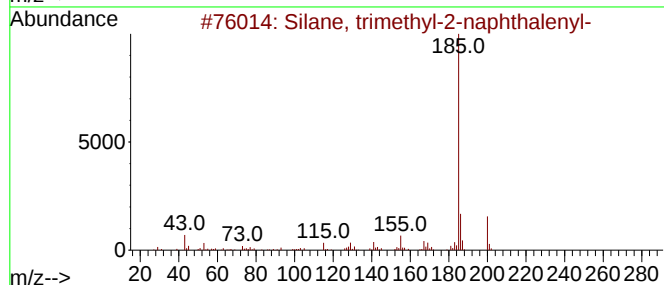
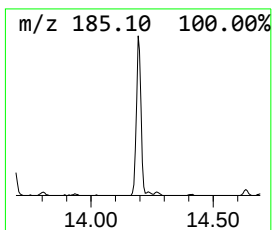
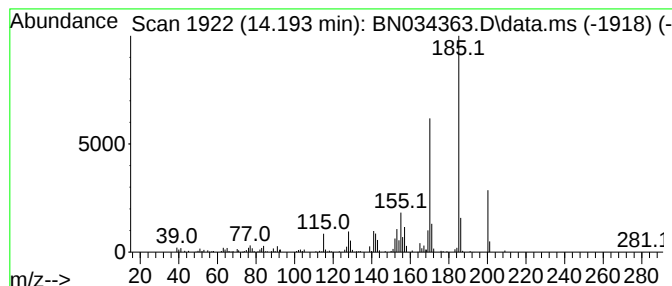
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Silane, trimethyl-2-naphtha... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.193	2.24 ng/ul	235602	Acenaphthene-d10	14.034

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Silane, trimethyl-2-naphthalenyl-	200	C13H16Si	018052-85-2	50
2			1H-Pyrazole, 5-(t-butyl)-3-phenyl-	200	C13H16N2	1000261-34-8	46
3			Benzaldehyde, 3-bromo-	184	C7H5BrO	003132-99-8	43
4			4'-(1-Pyrrolyl)acetophenone	185	C12H11NO	022106-37-2	38
5			4-Chlorothiopheno[2,3-b]pyridine-N...	185	C7H4ClNOS	025557-54-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100224\
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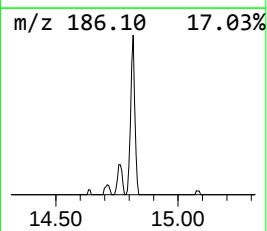
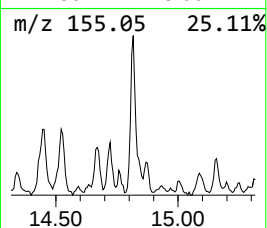
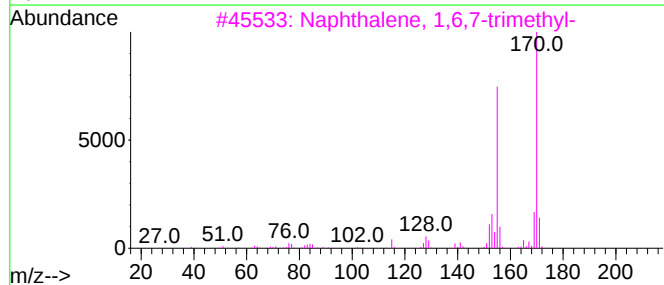
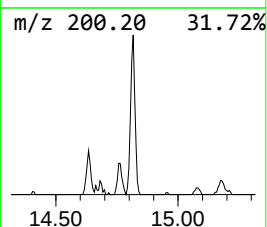
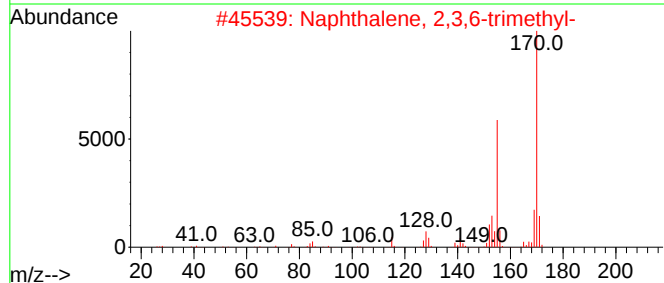
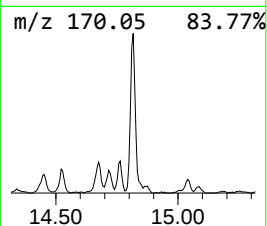
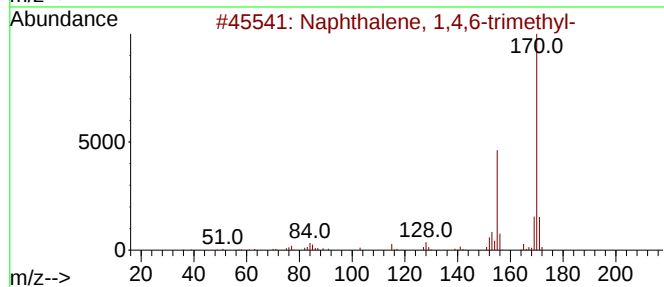
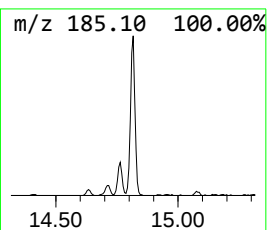
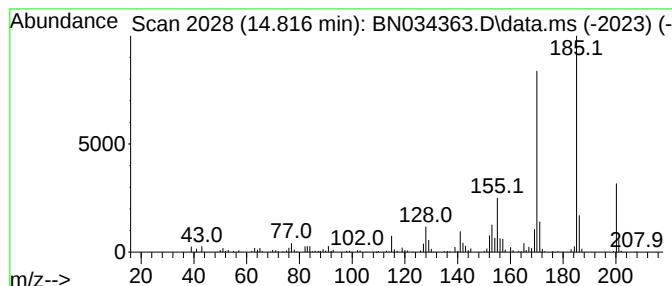
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Naphthalene, 1,4,6-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.816	2.69 ng/ul	282548	Acenaphthene-d10	14.034

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	50
2			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	50
3			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	46
4			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	46
5			4,6,8-Trimethylazulene	170	C13H14	000941-81-1	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100224\
Data File : BN034363.D
Acq On : 03 Oct 2024 09:58
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ALS Vial : 50 Sample Multiplier: 1

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ClientSampleId :
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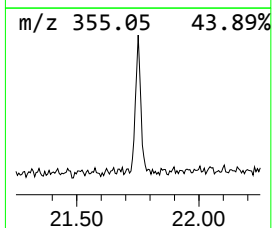
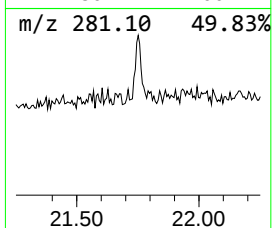
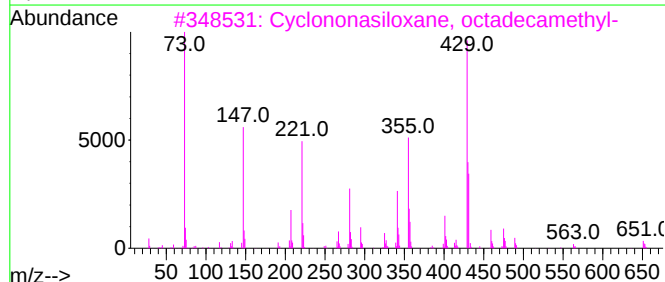
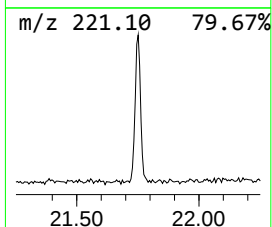
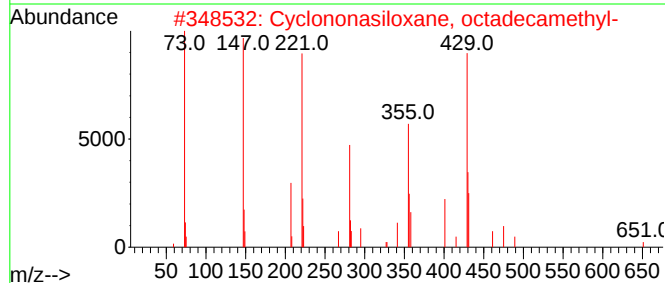
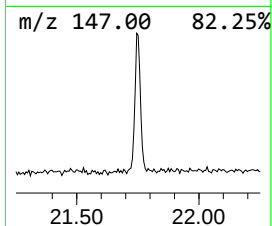
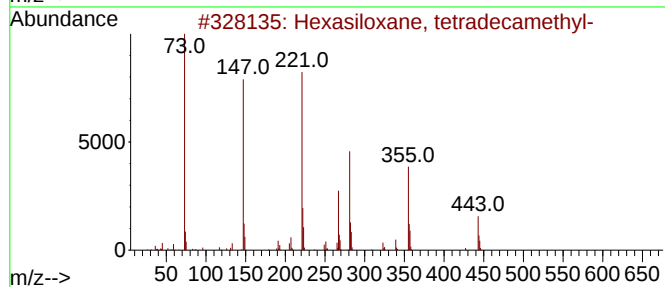
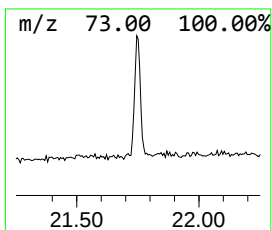
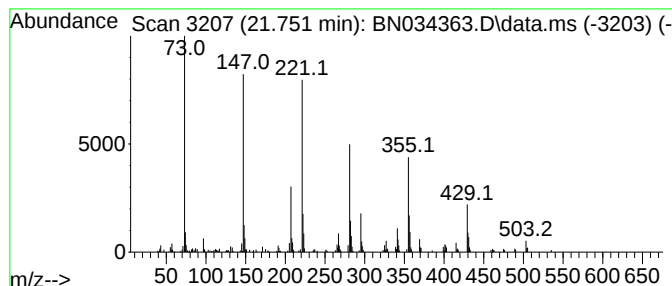
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Hexasiloxane, tetradecamethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.751	2.24 ng/ul	316008	Chrysene-d12	21.028

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexasiloxane, tetradecamethyl-	458	C14H42O5Si6	000107-52-8	83
2			Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	74
3			Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	64
4			1,1,1,5,7,7,7-Heptamethyl-3,3-bi...	444	C13H40O5Si6	038147-00-1	50
5			Mercaptoacetic acid, 2TMS deriva...	236	C8H20O2SSi2	006398-62-5	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100224\
Data File : BN034363.D
Acq On : 03 Oct 2024 09:58
Operator : RC/JU
Sample : P4018-04DL 20X
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DDCA2DL

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

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	5.446	3.0	ng/ul	186802	1	7.381	1259210	20.0
(DEL) Alkane: S...	6.564	3.2	ng/ul	199303	1	7.381	1259210	20.0
Butanoic acid, ...	6.922	3.0	ng/ul	188297	1	7.381	1259210	20.0
(DEL) Alkane: S...	7.028	10.0	ng/ul	630918	1	7.381	1259210	20.0
1-Hexanol, 2-et...	7.464	7.4	ng/ul	465225	1	7.381	1259210	20.0
(DEL) Alkane: S...	8.611	6.0	ng/ul	375877	1	7.381	1259210	20.0
(DEL) Alkane: S...	13.969	9.4	ng/ul	986147	3	14.034	2100970	20.0
Silane, trimeth...	14.193	2.2	ng/ul	235602	3	14.034	2100970	20.0
Naphthalene, 1,...	14.816	2.7	ng/ul	282548	3	14.034	2100970	20.0
Hexasiloxane, t...	21.751	2.2	ng/ul	316008	5	21.028	2825540	20.0