

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100224\
 Data File : BN034362.D
 Acq On : 03 Oct 2024 09:19
 Operator : RC/JU
 Sample : P4017-22
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DD8D8

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: BN034362.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.005	16	20	30	rVB	44192	63692	1.26%	0.113%
2	3.393	83	86	98	rBV	239174	372352	7.36%	0.662%
3	3.693	133	137	144	rVB	16717	25702	0.51%	0.046%
4	5.446	430	435	442	rVB4	14708	23129	0.46%	0.041%
5	5.758	483	488	496	rBV	28700	51631	1.02%	0.092%
6	6.493	610	613	617	rVB	21083	27878	0.55%	0.050%
7	6.581	617	628	643	rBV	628365	1079346	21.34%	1.918%
8	6.740	649	655	667	rBV	529017	817016	16.15%	1.452%
9	6.834	667	671	676	rVV	25676	39167	0.77%	0.070%
10	6.922	679	686	698	rVV	706055	1088505	21.52%	1.934%
11	7.028	698	704	712	rVV2	78085	146105	2.89%	0.260%
12	7.381	756	764	773	rBV	922646	1417112	28.02%	2.518%
13	7.463	773	778	790	rVB3	38146	89555	1.77%	0.159%
14	7.622	800	805	820	rVB	102177	201011	3.97%	0.357%
15	7.810	830	837	843	rBV	307359	467497	9.24%	0.831%
16	7.922	852	856	861	rVB3	23490	39003	0.77%	0.069%
17	8.004	861	870	875	rBV4	63830	144768	2.86%	0.257%
18	8.099	880	886	893	rBV	777761	1228648	24.29%	2.183%
19	8.375	929	933	939	rVB2	35132	53498	1.06%	0.095%
20	8.475	943	950	951	rVV3	17170	23658	0.47%	0.042%
21	8.522	951	958	968	rVV	584298	1013004	20.03%	1.800%
22	8.610	968	973	979	rVV	66935	105917	2.09%	0.188%
23	8.687	979	986	999	rVV	103680	225682	4.46%	0.401%
24	8.804	999	1006	1010	rVV7	13088	29264	0.58%	0.052%
25	8.869	1014	1017	1026	rVV8	14968	33400	0.66%	0.059%
26	8.957	1027	1032	1043	rVV6	23528	56843	1.12%	0.101%
27	9.110	1051	1058	1063	rVV2	67074	112587	2.23%	0.200%
28	9.163	1063	1067	1072	rVB3	13891	24414	0.48%	0.043%
29	9.240	1072	1080	1087	rBV	498841	844934	16.71%	1.501%
30	9.310	1087	1092	1097	rVV3	16305	30192	0.60%	0.054%
31	9.669	1149	1153	1163	rVV8	15185	44819	0.89%	0.080%
32	9.769	1163	1170	1179	rVV	772558	1288205	25.47%	2.289%
33	9.851	1179	1184	1192	rVB	38064	68214	1.35%	0.121%
34	10.040	1208	1216	1224	rBV2	48291	97908	1.94%	0.174%
35	10.140	1226	1233	1244	rBV	1237748	2373321	46.93%	4.217%
36	10.287	1251	1258	1270	rBV	653996	1209869	23.92%	2.150%

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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	10.540	1295	1301	1311	rVB	79920	147915	2.92%	0.263%
38	10.710	1326	1330	1336	rVV2	25830	40887	0.81%	0.073%
39	11.045	1382	1387	1390	rVV2	27110	49238	0.97%	0.087%
40	11.075	1390	1392	1405	rVV7	22361	42819	0.85%	0.076%
41	11.393	1440	1446	1449	rBV5	17911	27869	0.55%	0.050%
42	11.575	1472	1477	1483	rBV4	26551	52744	1.04%	0.094%
43	11.734	1501	1504	1509	rVB3	15472	25621	0.51%	0.046%
44	11.810	1512	1517	1522	rBV2	27302	43535	0.86%	0.077%
45	12.034	1547	1555	1562	rVB3	24927	66294	1.31%	0.118%
46	12.604	1645	1652	1658	rVV	40229	73437	1.45%	0.130%
47	12.792	1680	1684	1691	rVV9	12374	28252	0.56%	0.050%
48	12.875	1691	1698	1704	rVV2	29621	49923	0.99%	0.089%
49	13.087	1726	1734	1740	rVB8	21848	52030	1.03%	0.092%
50	13.210	1744	1755	1764	rBV8	19056	80020	1.58%	0.142%
51	13.351	1775	1779	1783	rVV3	22449	35707	0.71%	0.063%
52	13.404	1783	1788	1792	rVV6	18843	32480	0.64%	0.058%
53	13.463	1792	1798	1807	rVV	1263381	1742334	34.45%	3.096%
54	13.592	1816	1820	1824	rBV4	17812	27833	0.55%	0.049%
55	13.722	1835	1842	1852	rBV	1563072	2291667	45.31%	4.072%
56	13.969	1878	1884	1888	rBV	28863	44085	0.87%	0.078%
57	14.034	1888	1895	1910	rVV	1590141	2372649	46.91%	4.216%
58	14.145	1910	1914	1919	rVB	25954	36565	0.72%	0.065%
59	14.269	1927	1935	1946	rBV	350819	668259	13.21%	1.187%
60	14.375	1949	1953	1957	rVB4	19764	34699	0.69%	0.062%
61	14.575	1983	1987	1989	rBV	28741	46085	0.91%	0.082%
62	14.675	1996	2004	2010	rBV	262004	390918	7.73%	0.695%
63	14.716	2010	2011	2017	rVB6	17744	21405	0.42%	0.038%
64	14.934	2041	2048	2052	rVB	30558	43711	0.86%	0.078%
65	15.039	2055	2066	2079	rBV	1948484	2788304	55.13%	4.954%
66	15.169	2081	2088	2093	rBV	452744	644308	12.74%	1.145%
67	15.345	2111	2118	2122	rVB8	11760	29600	0.59%	0.053%
68	15.422	2122	2131	2139	rBV	1504045	2071073	40.95%	3.680%
69	15.798	2192	2195	2198	rBV	30102	39171	0.77%	0.070%
70	15.875	2203	2208	2213	rVB	64834	90490	1.79%	0.161%
71	15.981	2220	2226	2231	rBV	262210	358530	7.09%	0.637%
72	16.451	2300	2306	2320	rBV	3677906	5057545	100.00%	8.986%
73	16.592	2327	2330	2336	rVB3	29093	43456	0.86%	0.077%
74	16.792	2359	2364	2375	rVV	1792063	2622368	51.85%	4.659%
75	16.892	2375	2381	2393	rVV	2108359	2853025	56.41%	5.069%
76	17.333	2452	2456	2461	rBV	27254	35374	0.70%	0.063%

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Integration Parameters: LSCINT.P

Integrator: RTE
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77	17.428	2467	2472	2479	rVB	147329	192924	3.81%	0.343%
78	17.727	2518	2523	2532	rBV	288107	484291	9.58%	0.860%
79	18.010	2565	2571	2572	rBV2	40153	58171	1.15%	0.103%
80	18.163	2594	2597	2604	rVB5	21389	35327	0.70%	0.063%
81	18.427	2637	2642	2648	rVB7	23507	42485	0.84%	0.075%
82	18.563	2661	2665	2671	rVB2	60879	98285	1.94%	0.175%
83	18.675	2681	2684	2691	rVB	28383	44492	0.88%	0.079%
84	18.763	2696	2699	2704	rVB2	27736	34833	0.69%	0.062%
85	18.833	2707	2711	2715	rBV2	31506	43008	0.85%	0.076%
86	18.874	2715	2718	2720	rBV3	20430	23322	0.46%	0.041%
87	19.027	2740	2744	2752	rBV	90425	145867	2.88%	0.259%
88	19.204	2767	2774	2781	rBV2	2480078	3361518	66.47%	5.973%
89	19.263	2781	2784	2787	rVV	40164	55530	1.10%	0.099%
90	19.298	2788	2790	2794	rVB3	28496	28089	0.56%	0.050%
91	19.804	2873	2876	2881	rVB	54763	60148	1.19%	0.107%
92	19.922	2894	2896	2902	rVB6	36382	40491	0.80%	0.072%
93	20.304	2958	2961	2965	rVB	65971	68334	1.35%	0.121%
94	20.763	3035	3039	3046	rBV2	300728	458051	9.06%	0.814%
95	21.027	3079	3084	3095	rBV	2174956	3023144	59.77%	5.371%
96	21.251	3119	3122	3127	rVB	79390	100230	1.98%	0.178%
97	21.774	3207	3211	3221	rVB3	90347	158719	3.14%	0.282%
98	22.510	3332	3336	3343	rVB	102070	168703	3.34%	0.300%
99	23.557	3506	3514	3529	rVB	1742785	3875591	76.63%	6.886%
100	23.739	3537	3545	3553	rBV	1601490	3519952	69.60%	6.254%

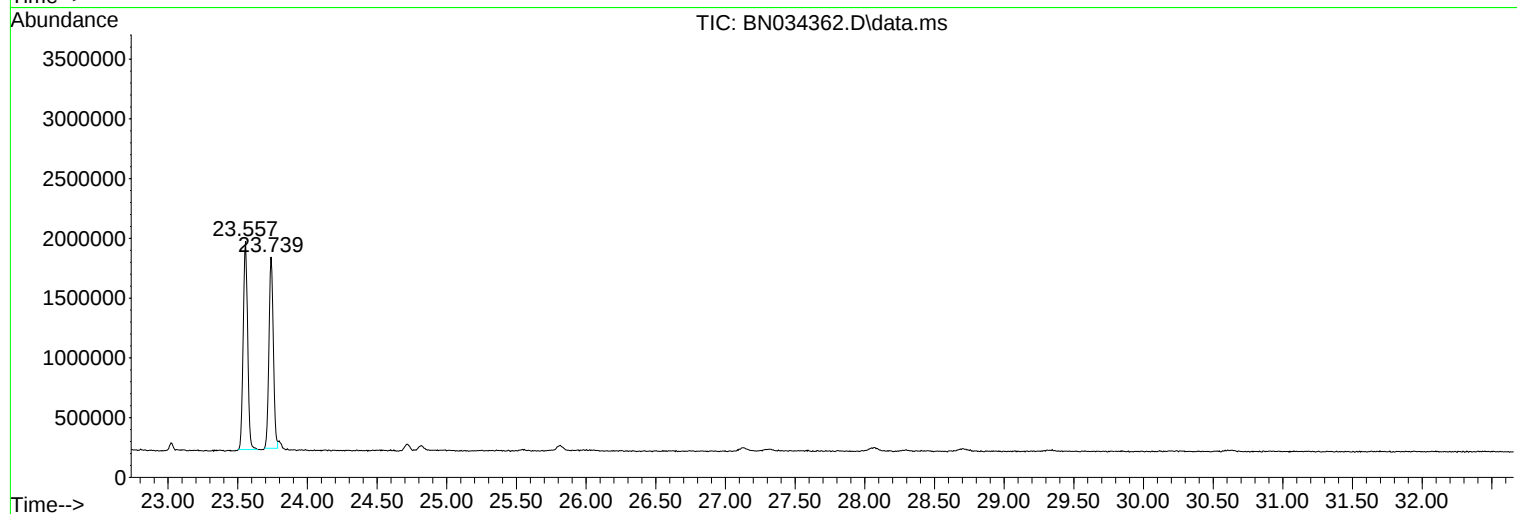
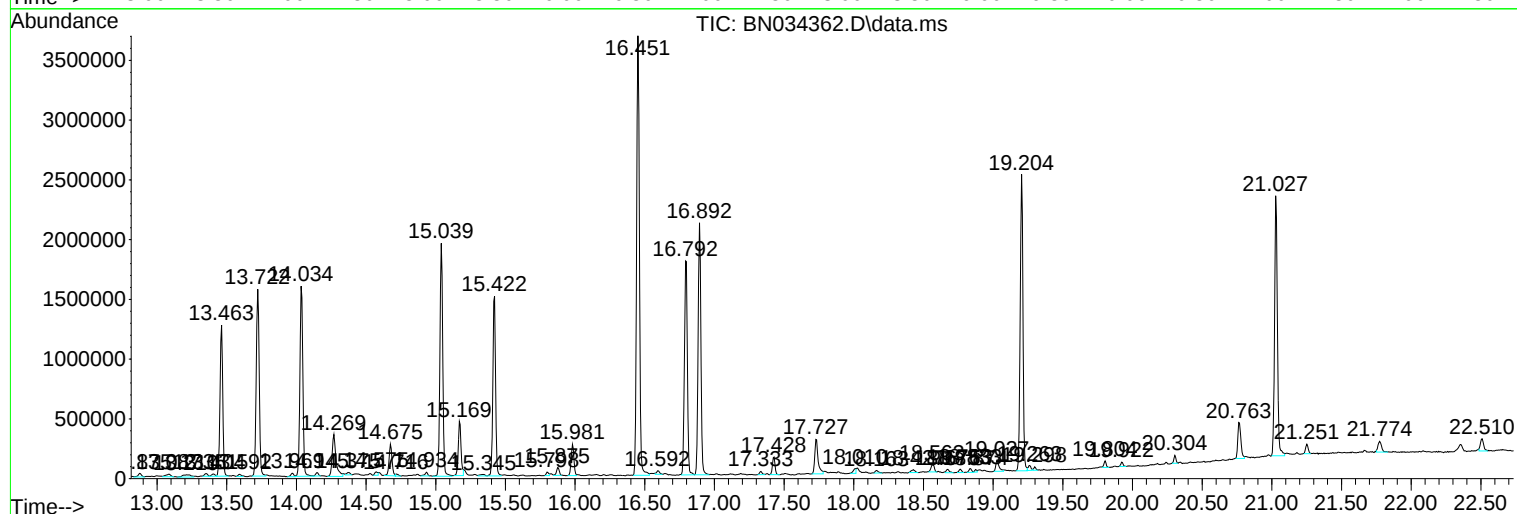
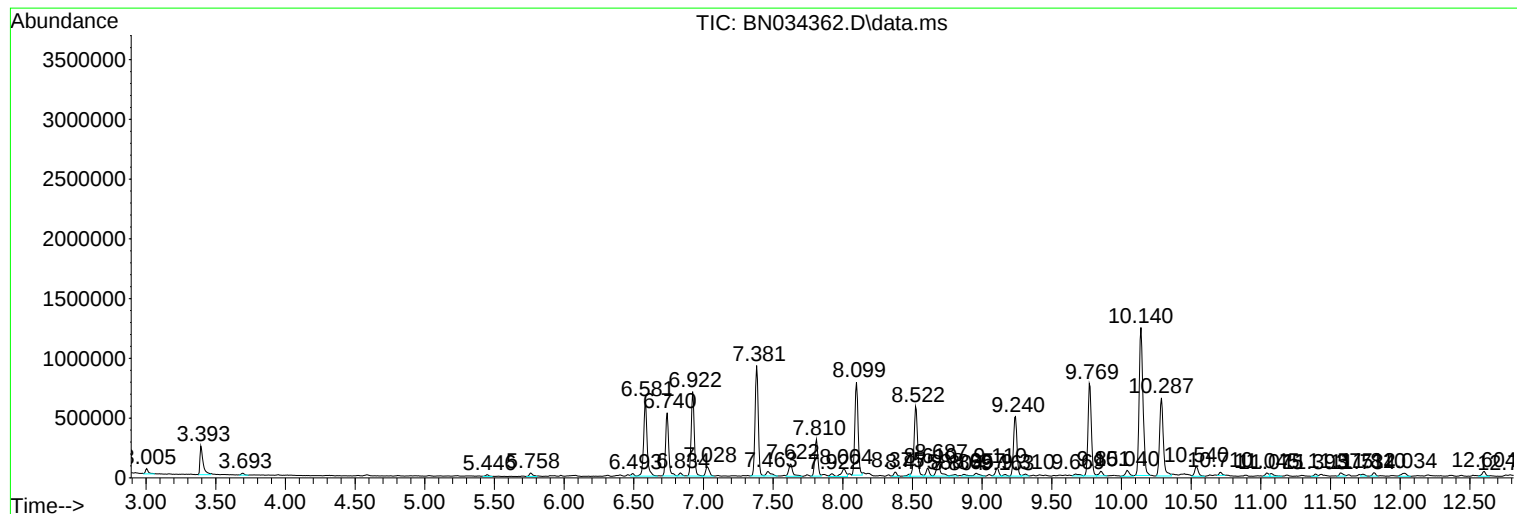
Sum of corrected areas: 56281576

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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



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100224\
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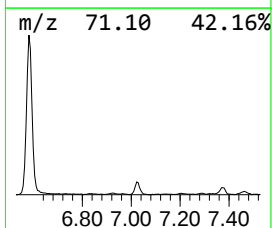
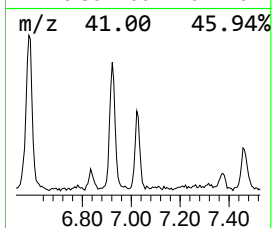
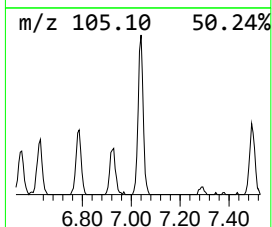
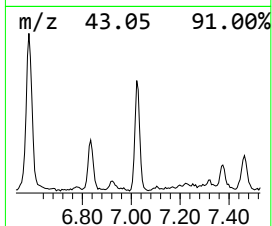
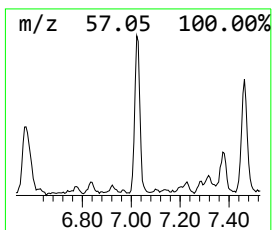
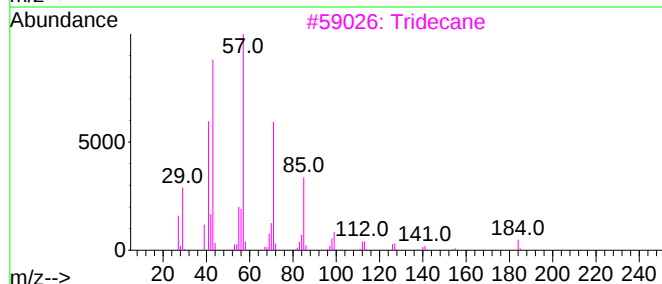
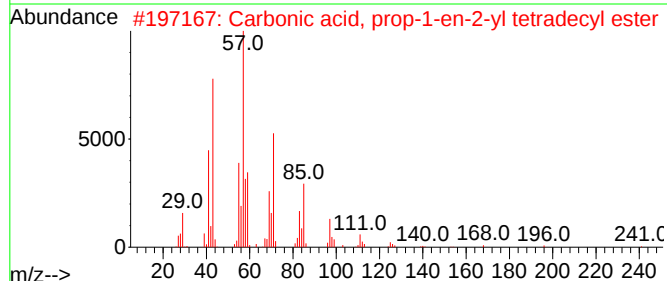
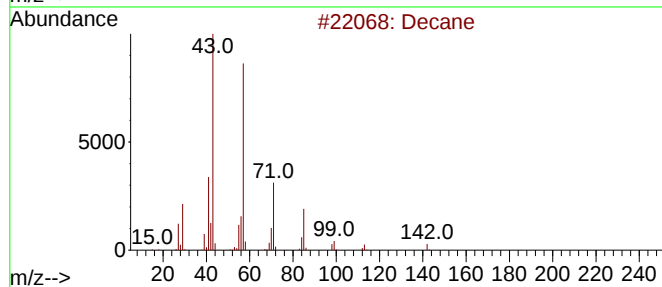
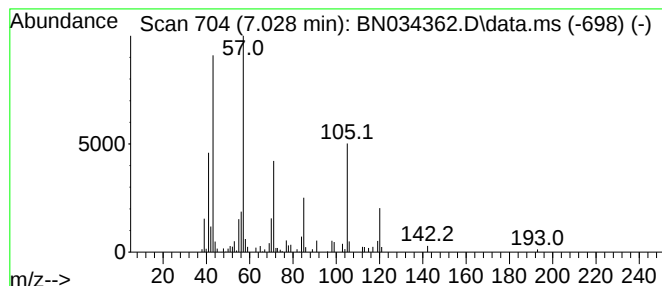
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.028	2.06 ng/ul	146105	1,4-Dichlorobenzene-d4	7.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane	142	C10H22	000124-18-5	64
2			Carbonic acid, prop-1-en-2-yl te...	298	C18H34O3	1000382-90-9	52
3			Tridecane	184	C13H28	000629-50-5	52
4			1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	50
5			Undecane	156	C11H24	001120-21-4	43



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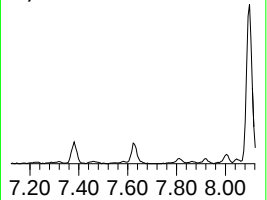
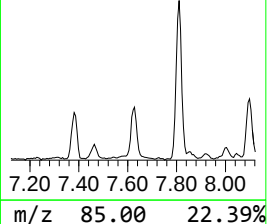
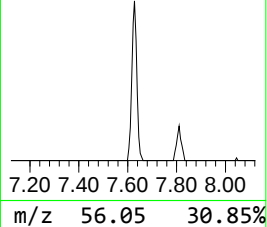
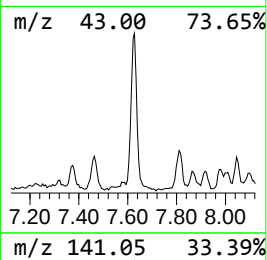
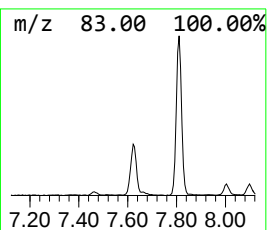
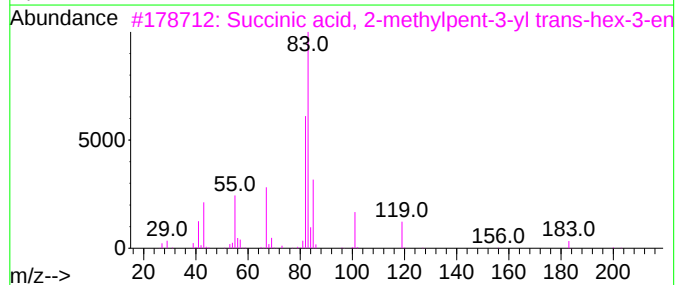
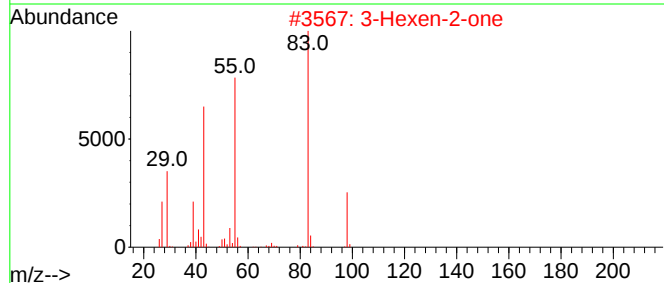
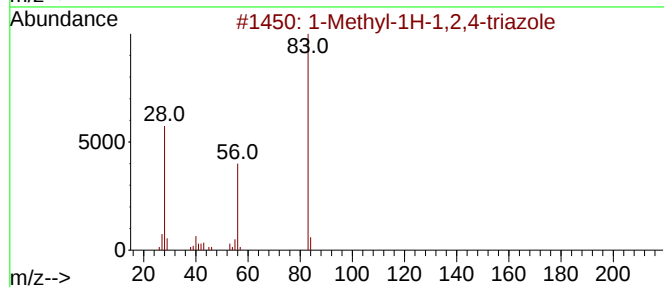
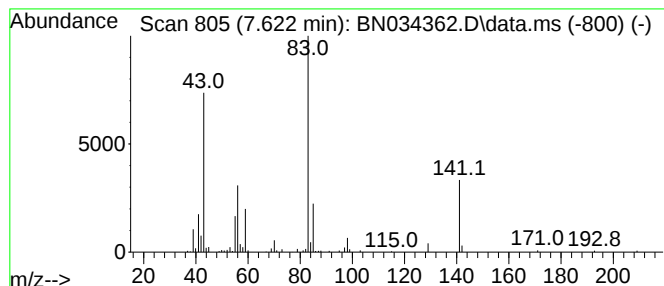
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.622	2.84 ng/ul	201011	1,4-Dichlorobenzene-d4	7.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Methyl-1H-1,2,4-triazole	83	C3H5N3	006086-21-1	37
2			3-Hexen-2-one	98	C6H10O	000763-93-9	25
3			Succinic acid, 2-methylpent-3-yl...	284	C16H28O4	1000391-11-2	23
4			2,4-Dimethylpentan-3-yl (E)-2-me...	198	C12H22O2	1000373-75-5	17
5			3-Methyl-2-butenic acid, pent-2...	164	C10H12O2	1000299-31-0	16



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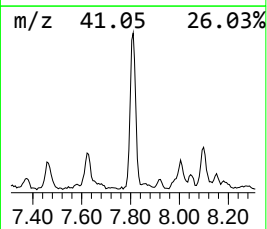
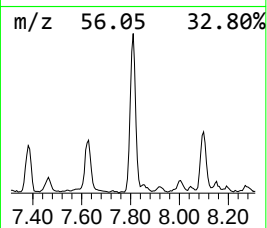
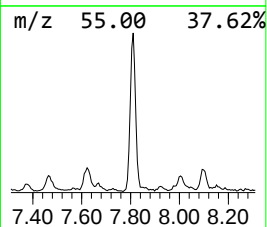
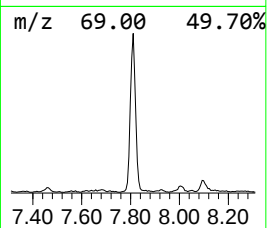
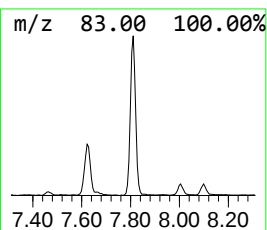
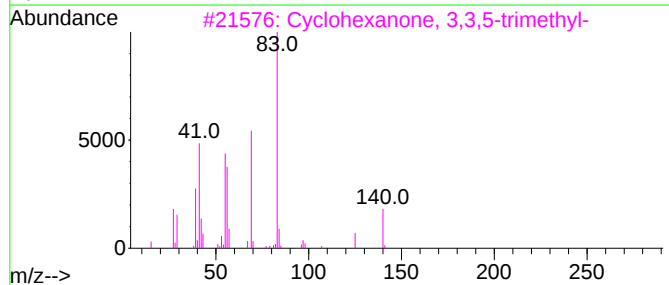
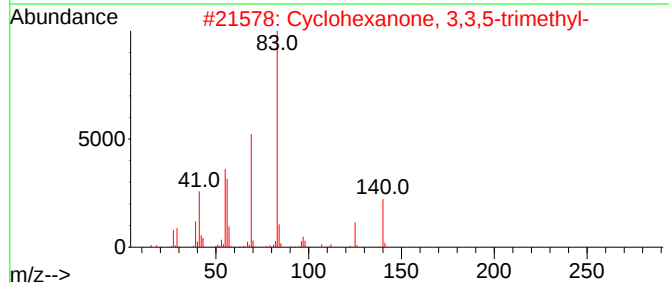
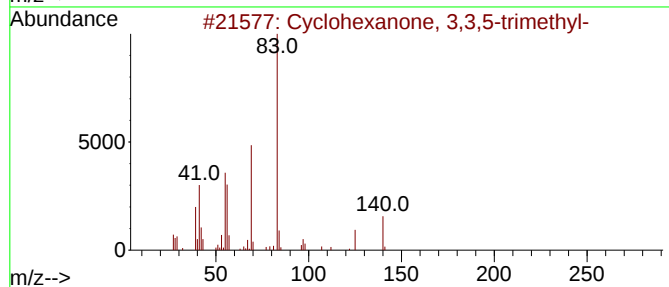
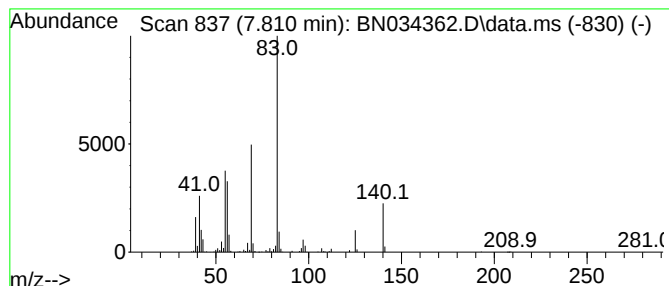
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN091124.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Cyclohexanone, 3,3,5-trimet... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.810	6.60 ng/ul	467497	1,4-Dichlorobenzene-d4	7.381

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	95
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			Cyclohexane, butyl-	140	C10H20	001678-93-9	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100224\
Data File : BN034362.D
Acq On : 03 Oct 2024 09:19
Operator : RC/JU
Sample : P4017-22
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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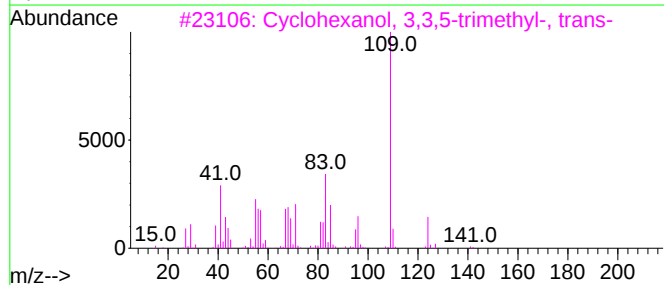
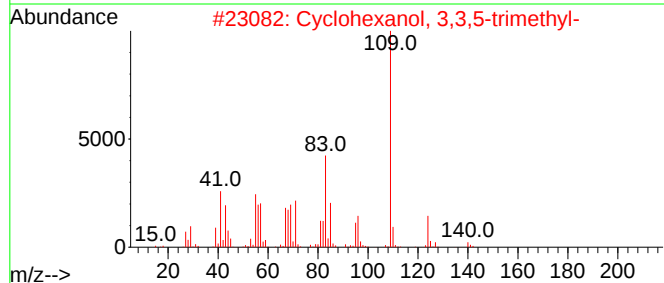
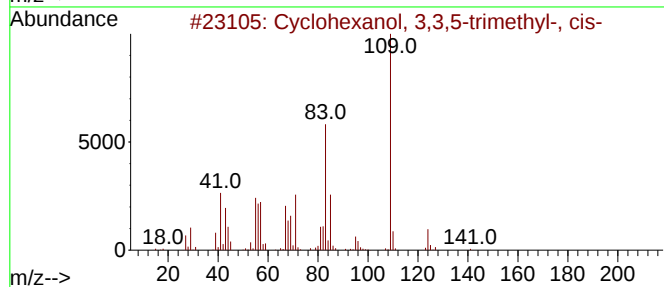
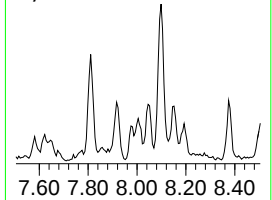
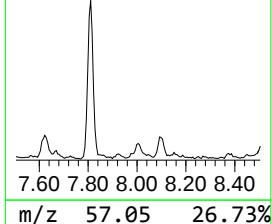
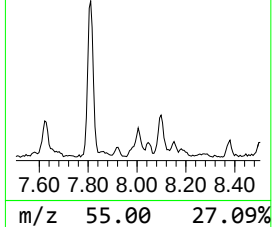
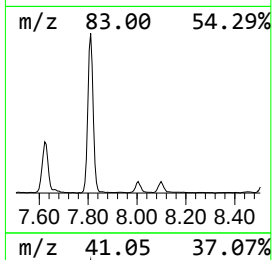
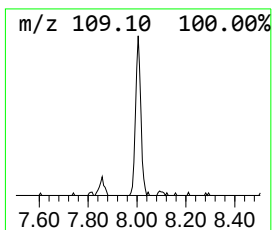
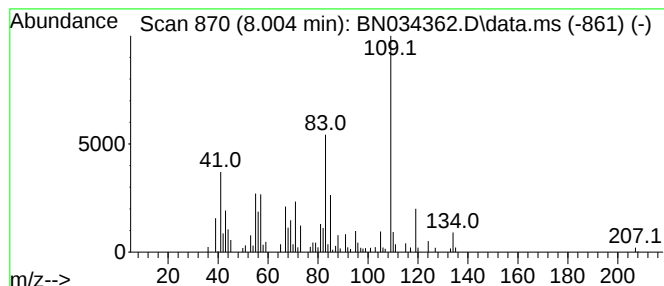
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Cyclohexanol, 3,3,5-trimeth... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.004	2.04 ng/ul	144768	1,4-Dichlorobenzene-d4	7.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanol, 3,3,5-trimethyl-, ...	142	C9H18O	000933-48-2	74
2			Cyclohexanol, 3,3,5-trimethyl-	142	C9H18O	000116-02-9	72
3			Cyclohexanol, 3,3,5-trimethyl-, ...	142	C9H18O	000767-54-4	59
4			Cyclopropanecarboxylic acid, 2,2...	168	C10H16O2	033383-56-1	37
5			2-Amino-4-methylpyrimidine	109	C5H7N3	000108-52-1	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100224\
Data File : BN034362.D
Acq On : 03 Oct 2024 09:19
Operator : RC/JU
Sample : P4017-22
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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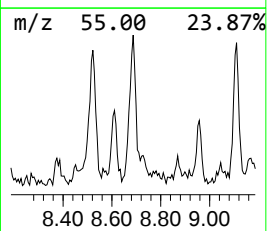
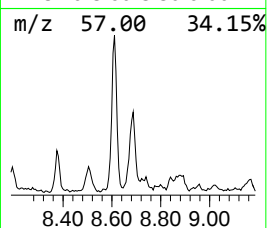
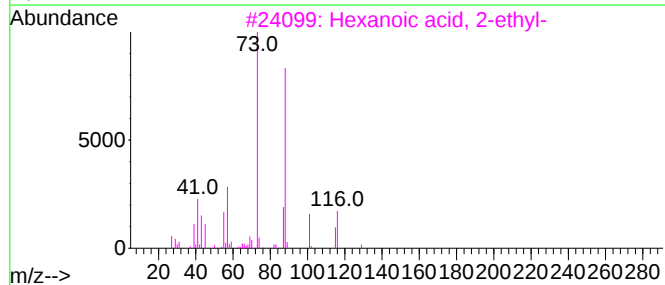
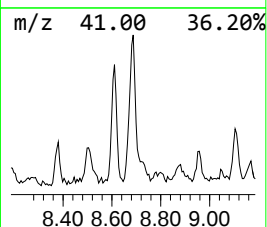
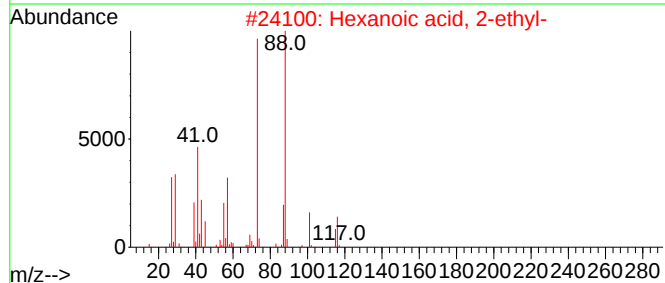
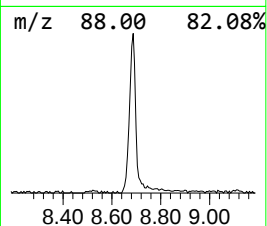
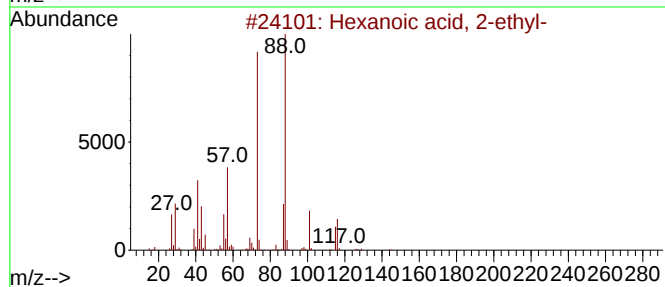
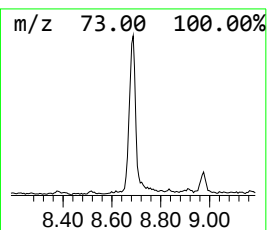
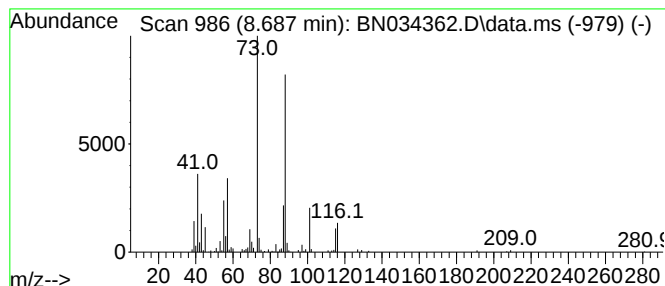
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Hexanoic acid, 2-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.687	3.19 ng/ul	225682	1,4-Dichlorobenzene-d4	7.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	90
2			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	80
3			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	78
4			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	64
5			2-Ethyl-4-methylpentanoic acid	144	C8H16O2	000108-81-6	56



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100224\
Data File : BN034362.D
Acq On : 03 Oct 2024 09:19
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Sample : P4017-22
Misc :
ALS Vial : 32 Sample Multiplier: 1

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ClientSampleId :
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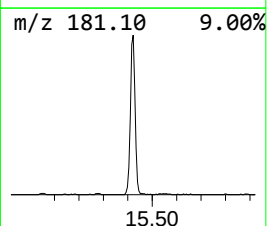
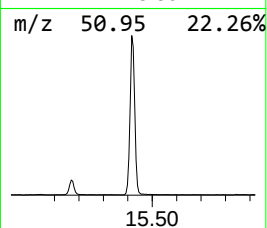
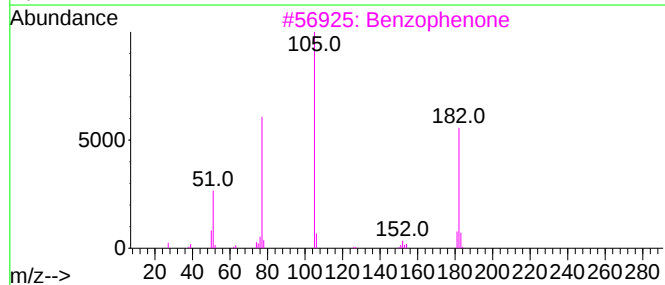
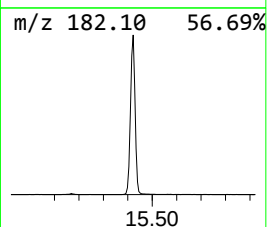
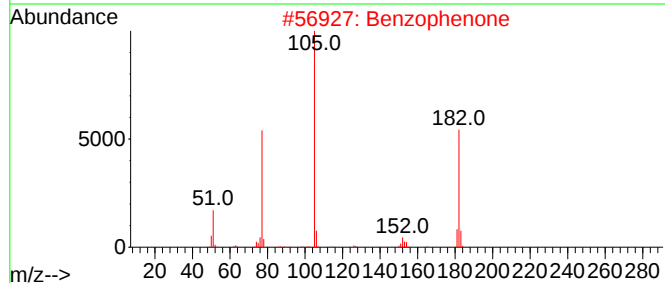
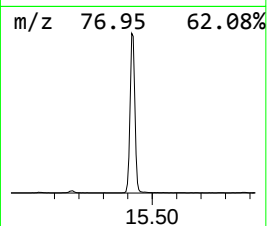
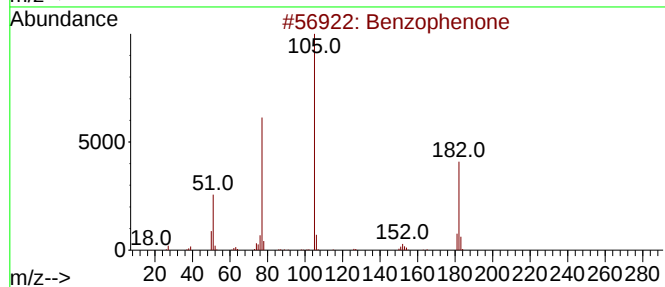
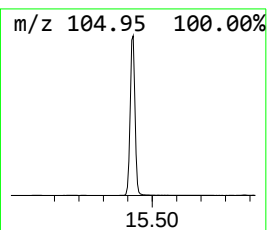
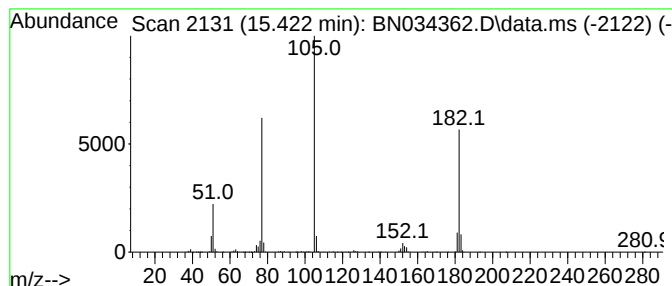
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzophenone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.422	15.80 ng/ul	2071070	Phenanthrene-d10	16.798

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	96
4			Benzophenone	182	C13H10O	000119-61-9	93
5			Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100224\
Data File : BN034362.D
Acq On : 03 Oct 2024 09:19
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Sample : P4017-22
Misc :
ALS Vial : 32 Sample Multiplier: 1

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ClientSampleId :
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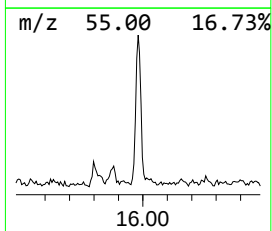
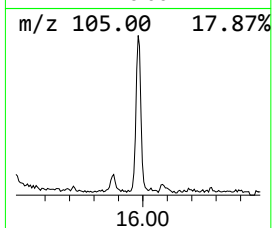
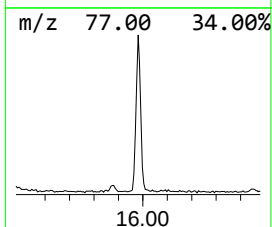
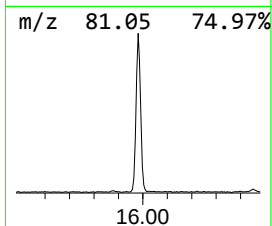
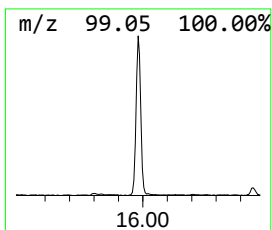
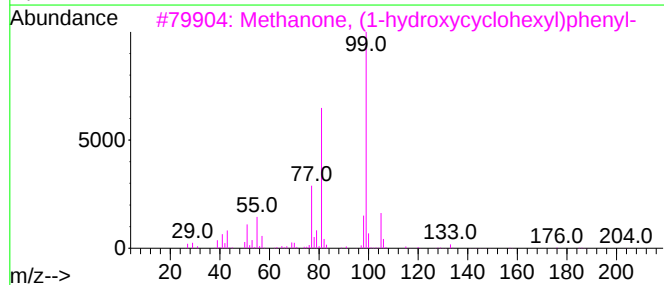
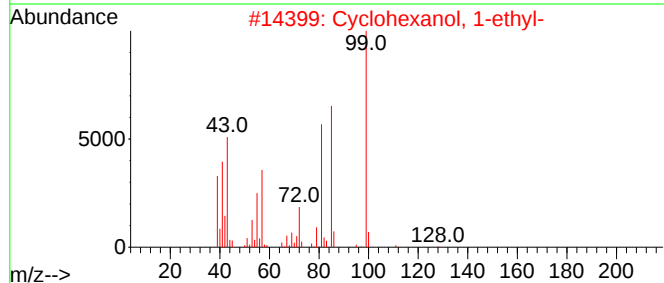
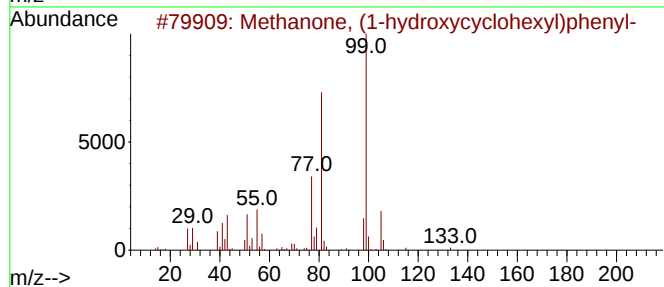
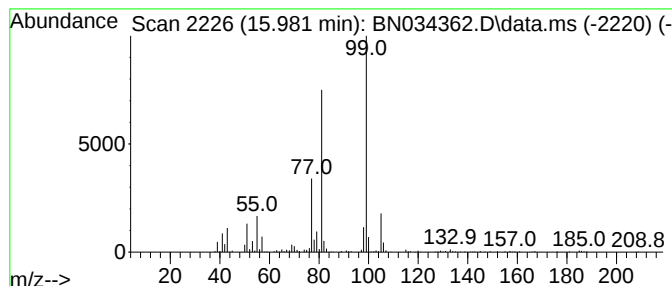
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Methanone, (1-hydroxycyclohexyl) Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.980	2.73 ng/ul	358530	Phenanthrene-d10	16.798

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methanone, (1-hydroxycyclohexyl)...	204	C13H16O2	000947-19-3	91
2			Cyclohexanol, 1-ethyl-	128	C8H16O	001940-18-7	59
3			Methanone, (1-hydroxycyclohexyl)...	204	C13H16O2	000947-19-3	52
4			1-Hydroxycyclohexanecarboxylic acid	144	C7H12O3	001123-28-0	45
5			3-Methylcyclohexyl methylphospho...	194	C8H16FO2P	113548-86-0	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100224\
Data File : BN034362.D
Acq On : 03 Oct 2024 09:19
Operator : RC/JU
Sample : P4017-22
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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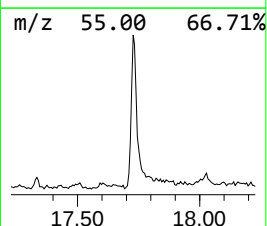
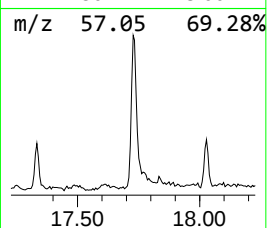
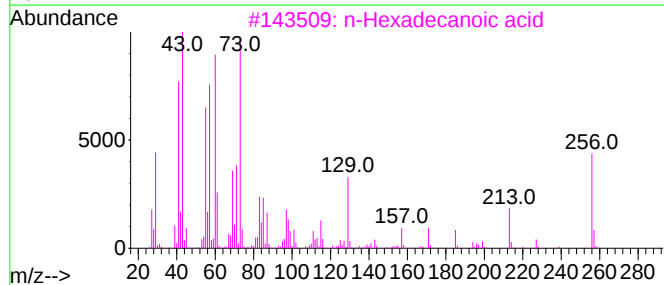
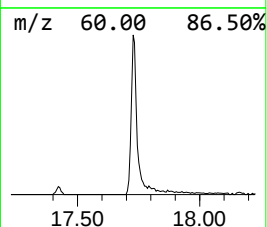
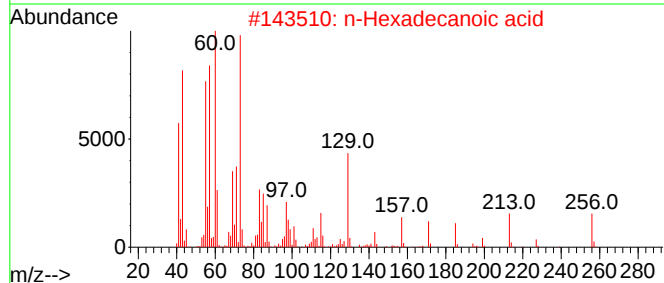
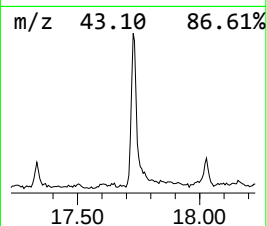
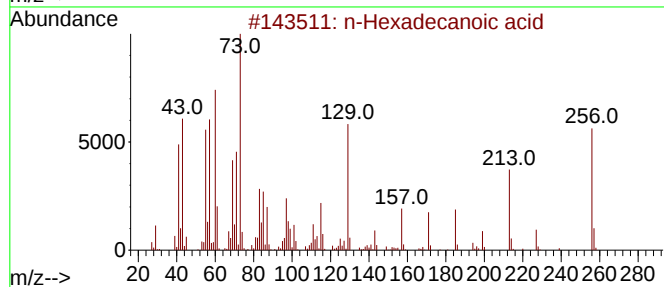
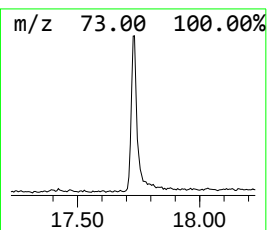
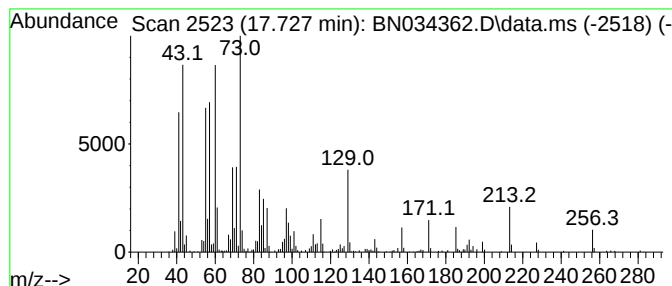
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.727	3.69 ng/ul	484291	Phenanthrene-d10	16.798

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			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
5			Tridecanoic acid	214	C13H26O2	000638-53-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100224\
Data File : BN034362.D
Acq On : 03 Oct 2024 09:19
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Sample : P4017-22
Misc :
ALS Vial : 32 Sample Multiplier: 1

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ClientSampleId :
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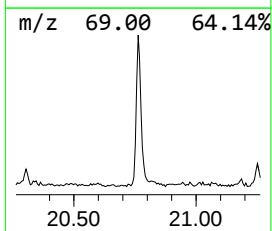
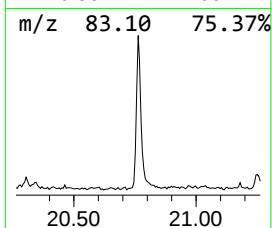
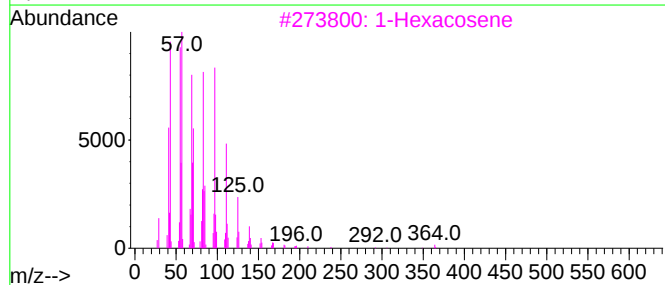
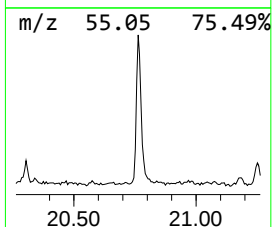
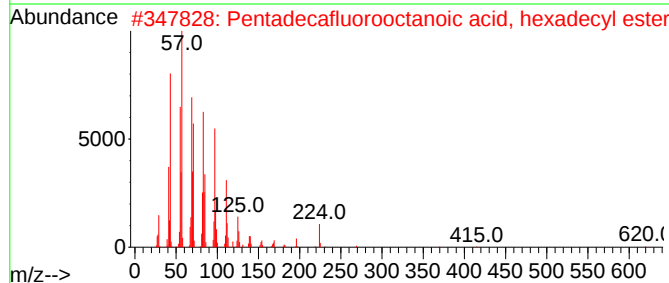
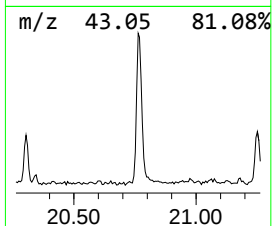
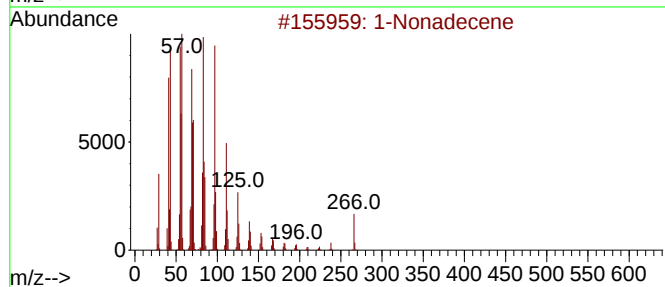
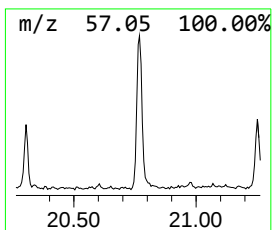
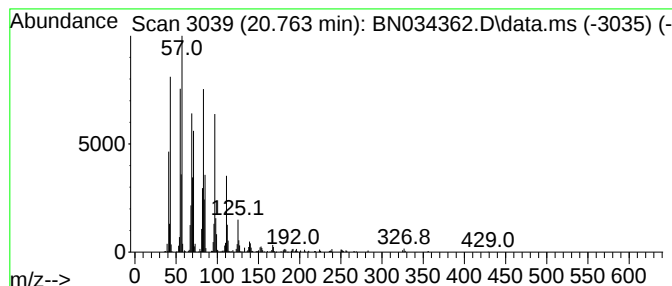
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.763	3.03 ng/ul	458051	Chrysene-d12	21.027

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Nonadecene	266	C19H38	018435-45-5	95
2			Pentadecafluorooctanoic acid, he...	638	C24H33F15O2	1000406-04-6	94
3			1-Hexacosene	364	C26H52	018835-33-1	91
4			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	91
5			Cyclohexadecane	224	C16H32	000295-65-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100224\
Data File : BN034362.D
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Sample : P4017-22
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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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...	7.028	2.1	ng/ul	146105	1	7.381	1417110	20.0
unknown-01	7.622	2.8	ng/ul	201011	1	7.381	1417110	20.0
Cyclohexanone, ...	7.810	6.6	ng/ul	467497	1	7.381	1417110	20.0
Cyclohexanol, 3...	8.004	2.0	ng/ul	144768	1	7.381	1417110	20.0
Hexanoic acid, ...	8.687	3.2	ng/ul	225682	1	7.381	1417110	20.0
Benzophenone	15.422	15.8	ng/ul	2071070	4	16.798	2622370	20.0
Methanone, (1-h...	15.980	2.7	ng/ul	358530	4	16.798	2622370	20.0
n-Hexadecanoic ...	17.727	3.7	ng/ul	484291	4	16.798	2622370	20.0
(DEL) Alkane: S...	20.763	3.0	ng/ul	458051	5	21.027	3023140	20.0