

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090424\
 Data File : BP021843.D
 Acq On : 05 Sep 2024 15:49
 Operator : RC/JU
 Sample : P3809-11
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YE3E5

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_P\Methods\SFAM-EPA-BP082324.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP021843.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.264	43	47	50	rBV	55810	71056	1.51%	0.123%
2	3.299	50	53	60	rVB	148504	179369	3.82%	0.311%
3	3.534	88	93	104	rBV	228971	340080	7.25%	0.590%
4	3.681	114	118	131	rBV	341625	497883	10.61%	0.864%
5	4.005	169	173	176	rVV	12947	18421	0.39%	0.032%
6	4.040	176	179	184	rVB7	5057	8828	0.19%	0.015%
7	4.452	244	249	254	rVB5	5339	11402	0.24%	0.020%
8	4.670	282	286	291	rBV3	6880	10474	0.22%	0.018%
9	4.793	302	307	312	rBV3	10375	16878	0.36%	0.029%
10	4.905	321	326	333	rVB	13516	23658	0.50%	0.041%
11	5.246	379	384	391	rBV2	9669	16936	0.36%	0.029%
12	5.875	487	491	496	rVB7	6295	10592	0.23%	0.018%
13	6.046	514	520	526	rVB4	23935	40376	0.86%	0.070%
14	6.775	637	644	649	rBV5	8478	15069	0.32%	0.026%
15	6.940	665	672	684	rBV	266070	431650	9.20%	0.749%
16	7.099	693	699	710	rBV	501370	805529	17.17%	1.399%
17	7.305	726	734	741	rBV	581947	1004613	21.41%	1.744%
18	7.369	741	745	754	rVB	94289	158714	3.38%	0.276%
19	7.769	806	813	819	rBV	878977	1435644	30.60%	2.493%
20	7.834	819	824	826	rVV6	9773	19202	0.41%	0.033%
21	7.858	826	828	833	rVB5	11410	13907	0.30%	0.024%
22	8.128	866	874	881	rBV10	9822	22607	0.48%	0.039%
23	8.469	924	932	939	rBV	700241	1111590	23.69%	1.930%
24	8.816	985	991	996	rBV	28574	41182	0.88%	0.072%
25	8.910	1000	1007	1017	rBV	553550	923412	19.68%	1.603%
26	9.040	1025	1029	1032	rVB3	7392	9178	0.20%	0.016%
27	9.158	1045	1049	1056	rVB2	23438	36205	0.77%	0.063%
28	9.358	1079	1083	1093	rVB5	14121	27101	0.58%	0.047%
29	9.475	1096	1103	1112	rBV	54992	104211	2.22%	0.181%
30	9.634	1123	1130	1139	rBV	427026	717899	15.30%	1.246%
31	9.863	1165	1169	1177	rVB	6375	13004	0.28%	0.023%
32	10.175	1213	1222	1231	rBV	745796	1293749	27.57%	2.246%
33	10.552	1278	1286	1295	rBV	1212112	2071780	44.16%	3.597%
34	10.681	1300	1308	1322	rBV	817859	1536700	32.75%	2.668%
35	10.910	1344	1347	1352	rBV7	5659	9589	0.20%	0.017%
36	11.040	1364	1369	1371	rBV6	5803	10051	0.21%	0.017%

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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_P\Methods\SFAM-EPA-BP082324.MA.M
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37	11.087	1372	1377	1382	rBV2	45110	76487	1.63%	0.133%
38	11.169	1389	1391	1401	rVB7	9410	17983	0.38%	0.031%
39	11.293	1401	1412	1422	rBV	339462	731983	15.60%	1.271%
40	11.893	1510	1514	1518	rVB4	13114	19826	0.42%	0.034%
41	12.051	1538	1541	1543	rBV4	7409	8741	0.19%	0.015%
42	12.257	1574	1576	1584	rVB8	8279	16980	0.36%	0.029%
43	12.428	1600	1605	1610	rVB2	20772	34942	0.74%	0.061%
44	12.493	1610	1616	1625	rBV10	18047	39741	0.85%	0.069%
45	12.834	1670	1674	1676	rBV5	7338	11421	0.24%	0.020%
46	12.957	1691	1695	1699	rBV6	12022	18144	0.39%	0.032%
47	13.004	1701	1703	1710	rVV8	10305	19655	0.42%	0.034%
48	13.557	1793	1797	1798	rBV4	12788	17533	0.37%	0.030%
49	13.810	1834	1840	1849	rBV	1395036	1975161	42.10%	3.429%
50	14.093	1878	1888	1902	rBV	1354341	2262894	48.23%	3.929%
51	14.334	1923	1929	1935	rBV	243263	366001	7.80%	0.635%
52	14.404	1935	1941	1952	rBV	1671185	2737673	58.35%	4.753%
53	14.622	1972	1978	1979	rBV	399194	555222	11.83%	0.964%
54	14.869	2017	2020	2027	rVB3	45880	82292	1.75%	0.143%
55	15.416	2106	2113	2125	rBV2	1757974	3187541	67.94%	5.534%
56	15.534	2127	2133	2140	rBV	577450	948042	20.21%	1.646%
57	15.816	2178	2181	2185	rVB2	50661	61602	1.31%	0.107%
58	16.075	2222	2225	2227	rBV4	30050	43357	0.92%	0.075%
59	16.275	2256	2259	2262	rBV	44907	59485	1.27%	0.103%
60	16.334	2266	2269	2276	rVV2	99821	152295	3.25%	0.264%
61	16.398	2277	2280	2284	rVB2	60418	77129	1.64%	0.134%
62	16.610	2311	2316	2321	rVB5	94363	158854	3.39%	0.276%
63	16.669	2323	2326	2331	rBV2	70476	98771	2.11%	0.171%
64	17.210	2412	2418	2424	rVB	2315193	3705831	78.99%	6.434%
65	17.310	2429	2435	2446	rVB	2174648	3570595	76.10%	6.199%
66	17.592	2480	2483	2488	rVB	91101	130154	2.77%	0.226%
67	17.916	2534	2538	2547	rVB2	235982	432699	9.22%	0.751%
68	18.045	2556	2560	2567	rVB	286470	457471	9.75%	0.794%
69	18.151	2573	2578	2583	rBV	1450633	1955104	41.67%	3.395%
70	19.480	2800	2804	2815	rVB	474523	726105	15.48%	1.261%
71	19.686	2834	2839	2848	rBV	3162543	4506594	96.05%	7.824%
72	20.369	2953	2955	2959	rVB	153454	149137	3.18%	0.259%
73	21.333	3115	3119	3126	rVB	994050	1434571	30.58%	2.491%
74	21.669	3170	3176	3184	rBV2	2570145	4506668	96.05%	7.825%
75	24.845	3708	3716	3727	rVB	1742913	4491134	95.72%	7.798%
76	25.074	3747	3755	3768	rVB2	1559741	4691782	100.00%	8.146%

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Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Title : SVOA CALIBRATION

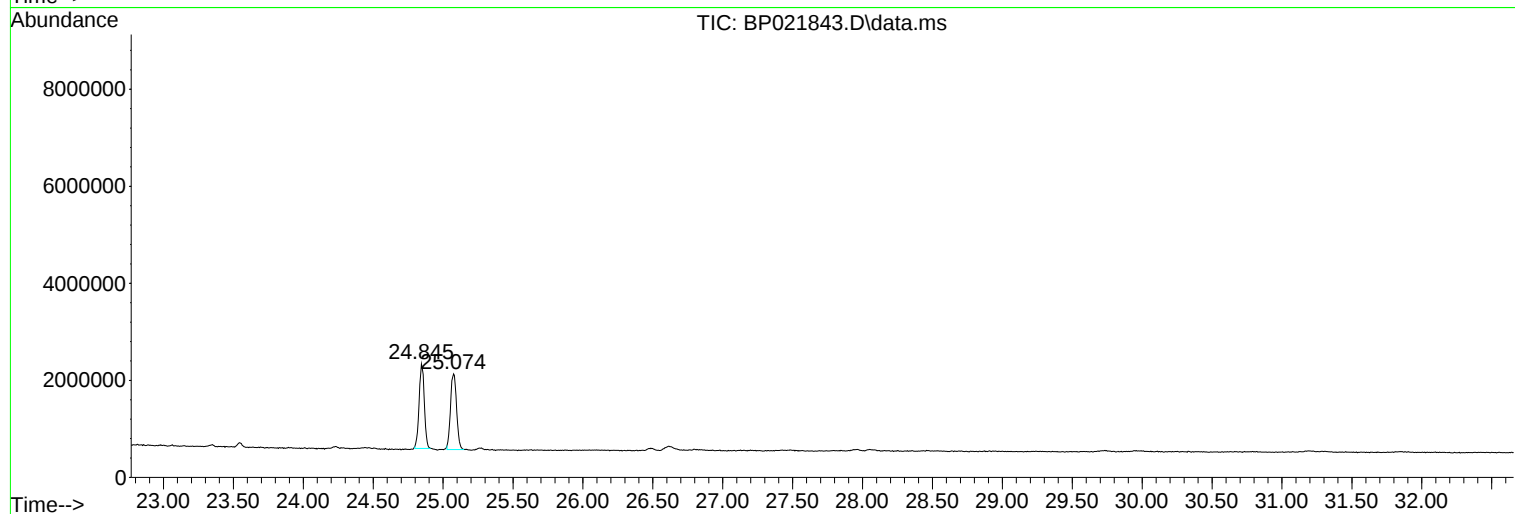
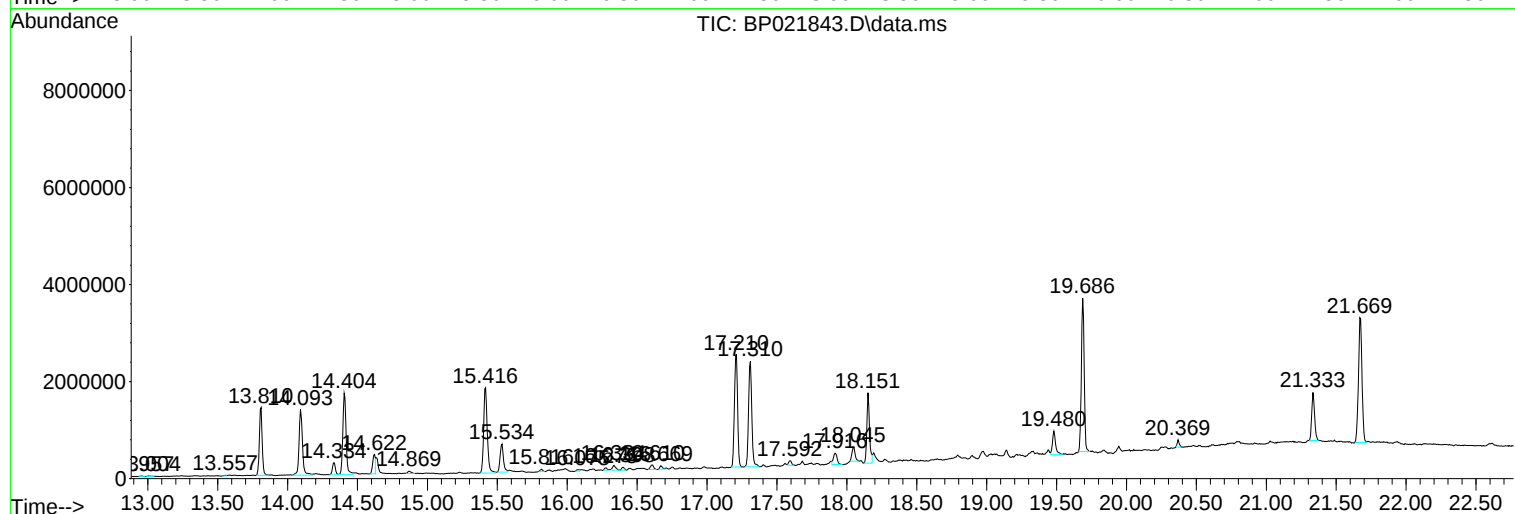
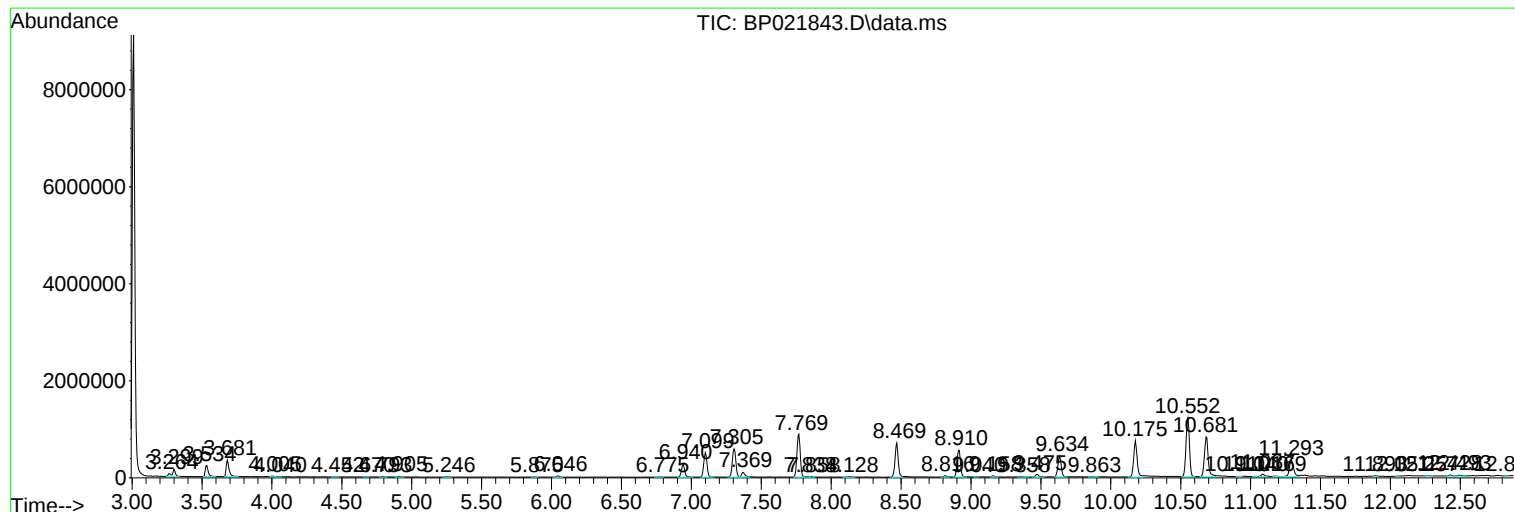
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090424\
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Acq On : 05 Sep 2024 15:49
Operator : RC/JU
Sample : P3809-11
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ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YE3E5

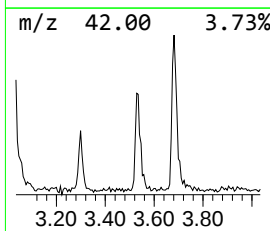
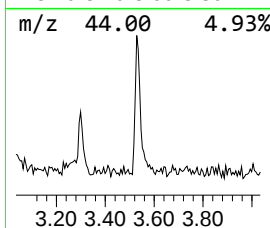
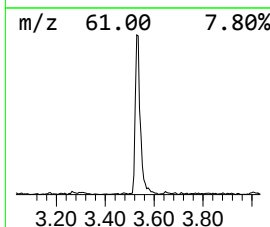
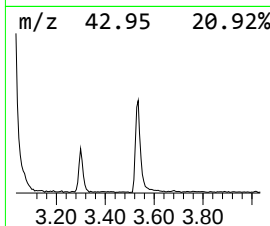
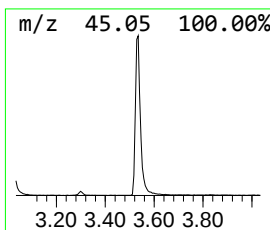
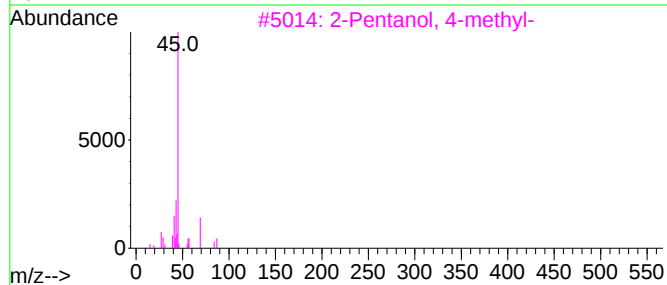
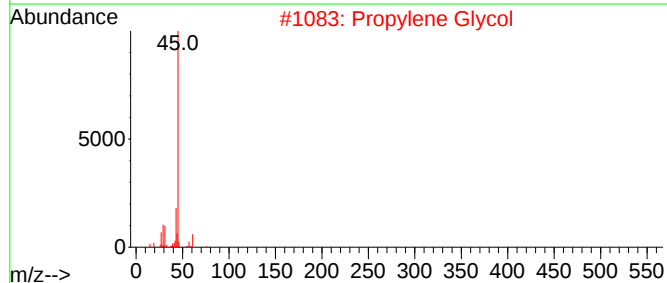
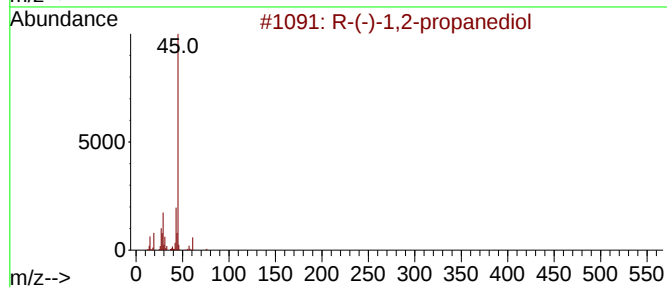
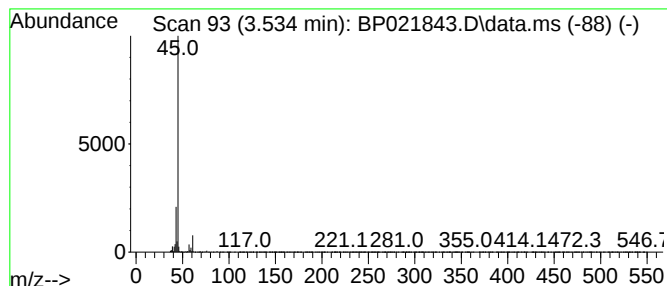
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 R-(-)-1,2-propanediol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.534	4.74 ng/ul	340080	1,4-Dichlorobenzene-d4	7.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	R-(-)-1,2-propanediol			76	C3H8O2	004254-14-2	72
2	Propylene Glycol			76	C3H8O2	000057-55-6	72
3	2-Pentanol, 4-methyl-			102	C6H14O	000108-11-2	40
4	Propanoic acid, 2-hydroxy-, meth...			104	C4H8O3	000547-64-8	40
5	Isopropyl Alcohol			60	C3H8O	000067-63-0	39



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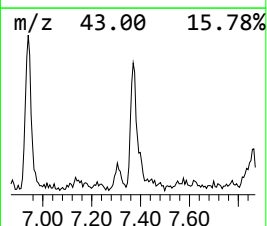
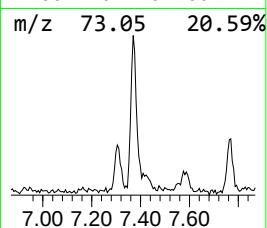
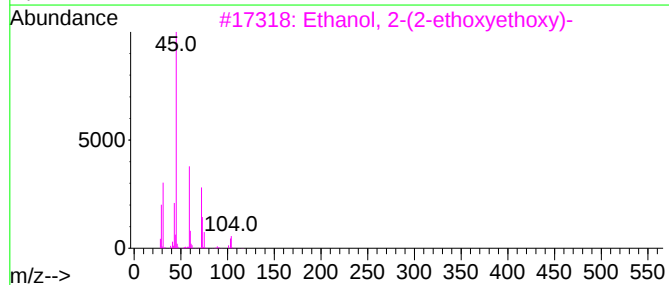
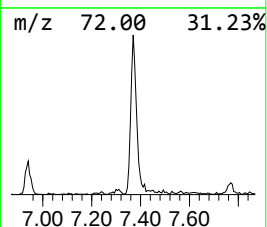
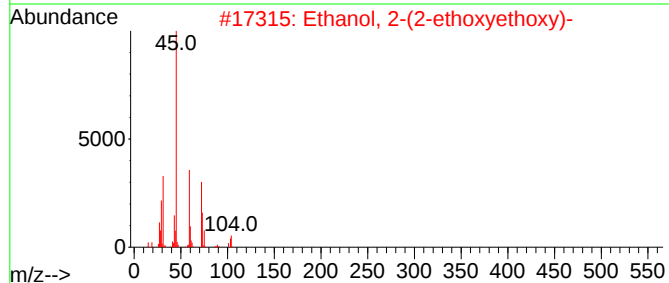
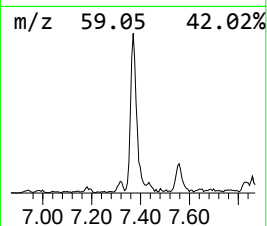
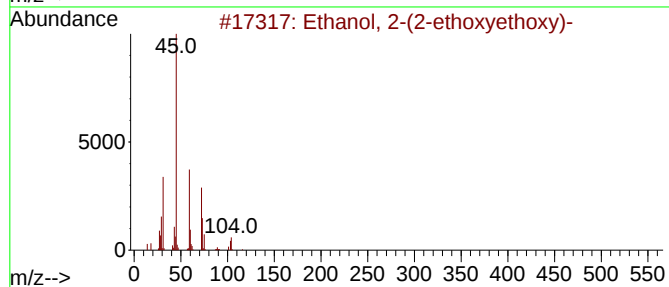
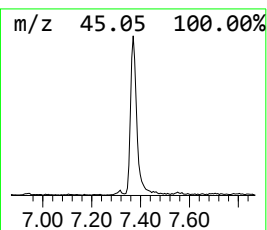
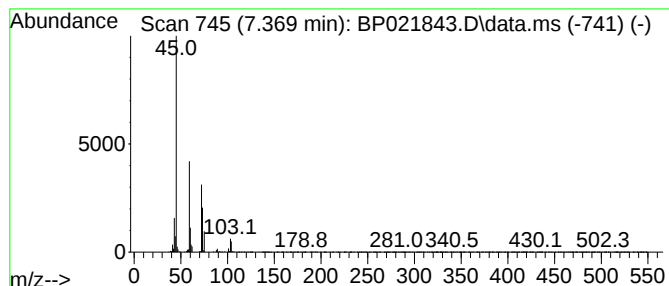
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.369	2.21 ng/ul	158714	1,4-Dichlorobenzene-d4	7.769

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	86
2			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	86
3			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	86
4			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	78
5			Diethyl carbitol	162	C8H18O3	000112-36-7	64



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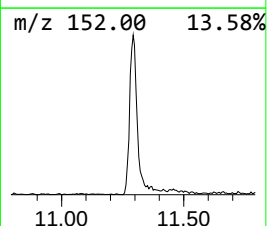
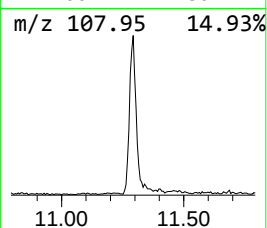
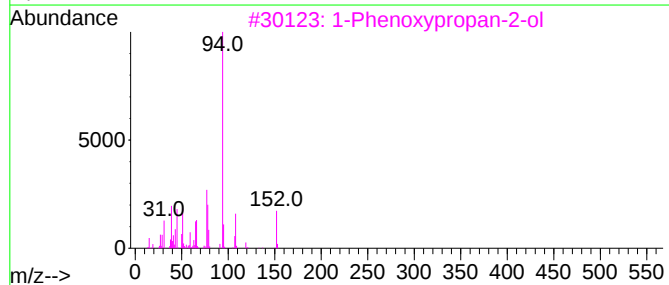
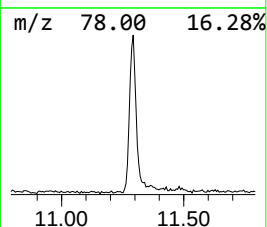
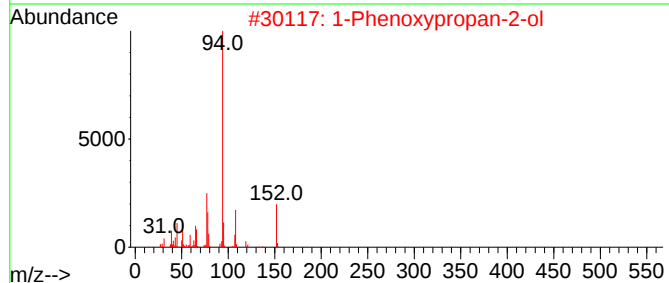
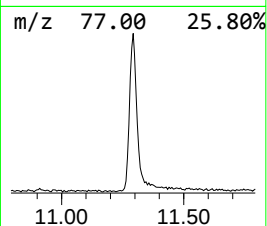
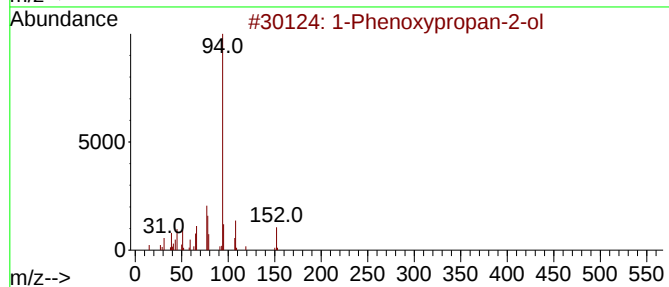
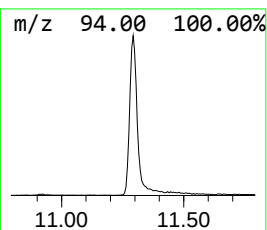
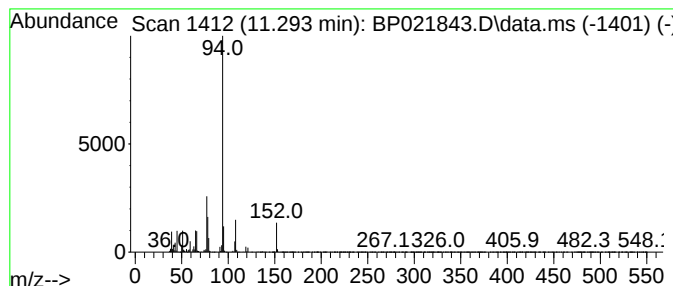
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.293	7.07 ng/ul	731983	Naphthalene-d8	10.552

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	95
4			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	91
5			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	64



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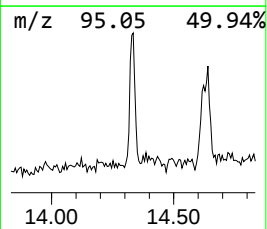
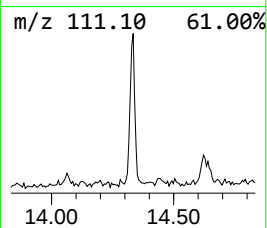
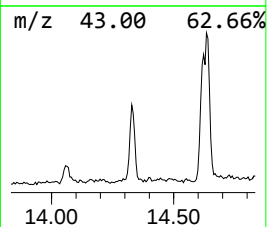
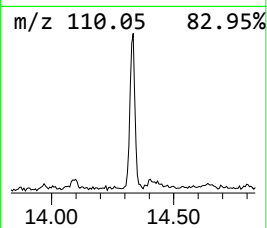
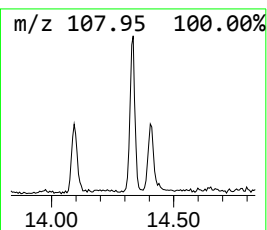
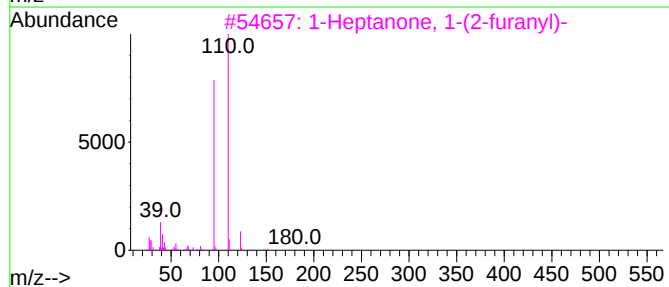
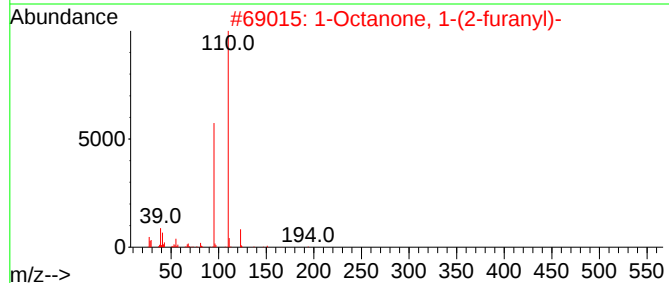
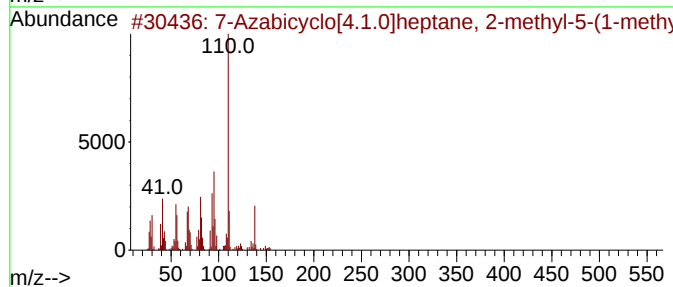
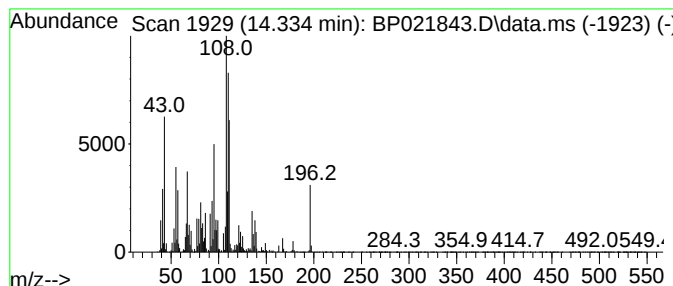
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.334	2.67 ng/ul	366001	Acenaphthene-d10	14.404

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7-Azabicyclo[4.1.0]heptane, 2-me...	153	C10H19N	055854-39-2	30
2			1-Octanone, 1-(2-furanyl)-	194	C12H18O2	005456-77-9	27
3			1-Heptanone, 1-(2-furanyl)-	180	C11H16O2	005466-40-0	27
4			Ethanone, 2-cyclopentyl-1-(1H-im...	178	C10H14N2O	069393-25-5	27
5			.alpha.-Campholenal	152	C10H16O	004501-58-0	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090424\
Data File : BP021843.D
Acq On : 05 Sep 2024 15:49
Operator : RC/JU
Sample : P3809-11
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YE3E5

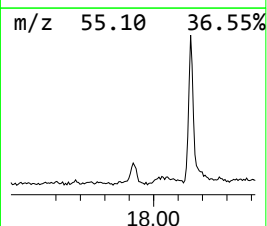
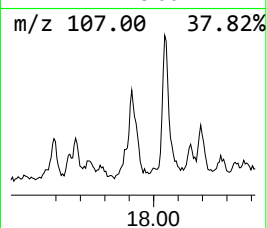
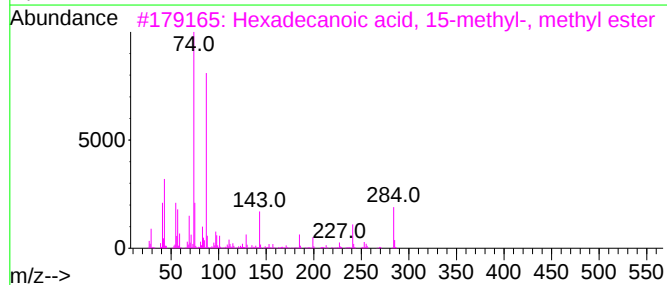
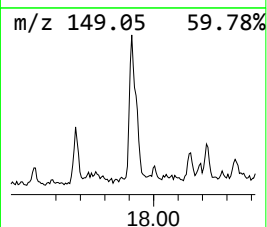
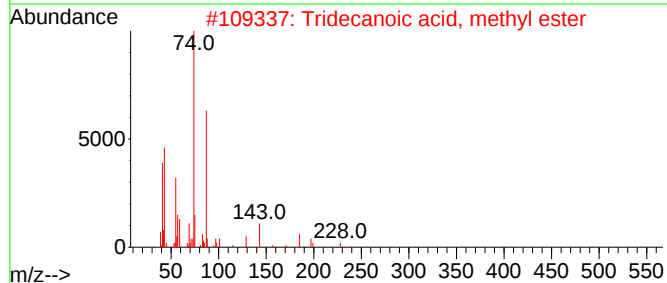
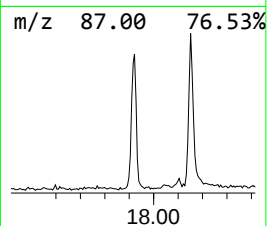
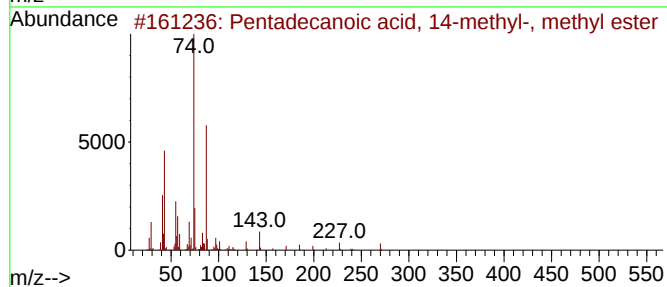
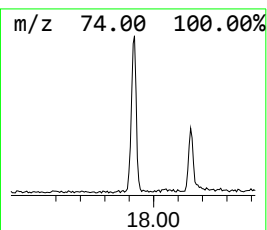
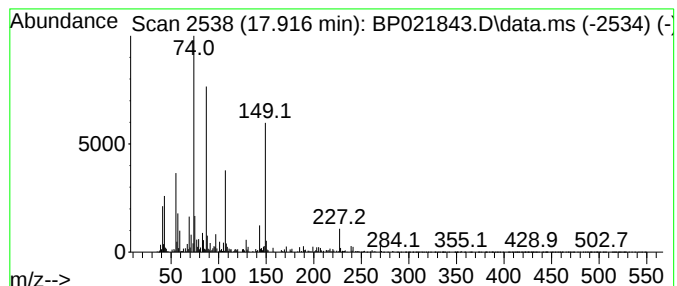
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Pentadecanoic acid, 14-meth... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.916	2.34 ng/ul	432699	Phenanthrene-d10	17.210

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	95
2			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	83
3			Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	83
4			Methyl tetradecanoate	242	C15H30O2	000124-10-7	62
5			Methyl tetradecanoate	242	C15H30O2	000124-10-7	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090424\
Data File : BP021843.D
Acq On : 05 Sep 2024 15:49
Operator : RC/JU
Sample : P3809-11
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YE3E5

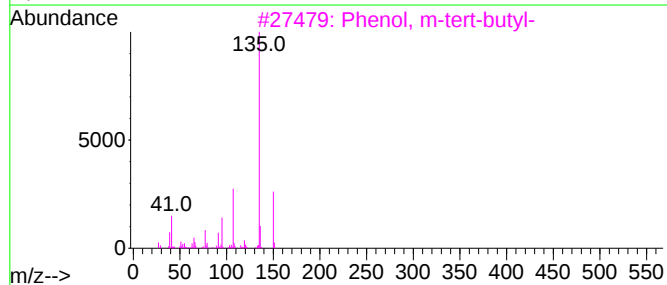
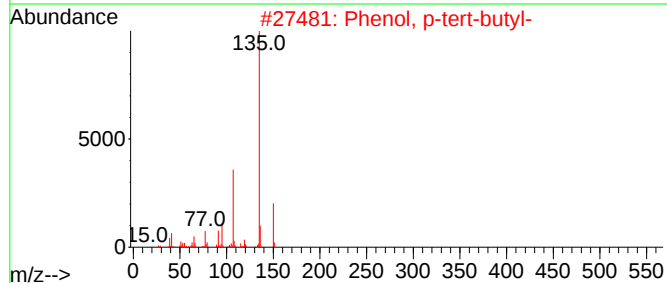
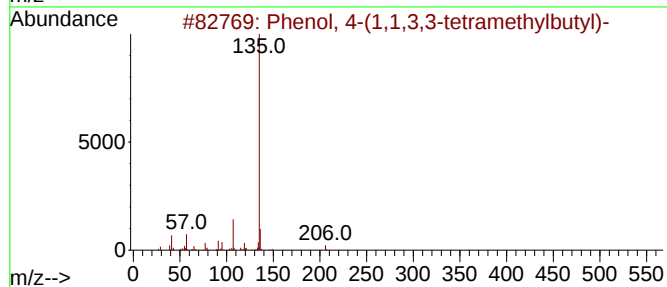
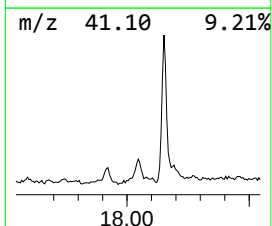
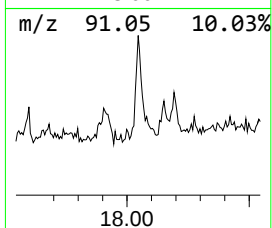
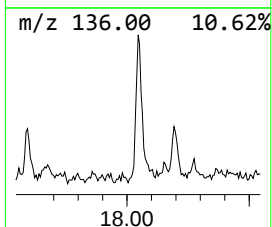
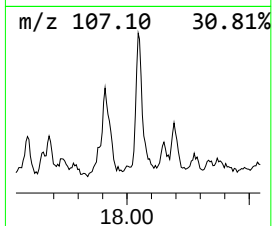
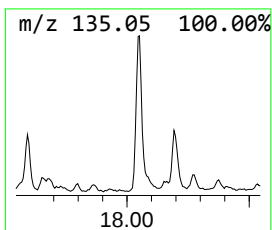
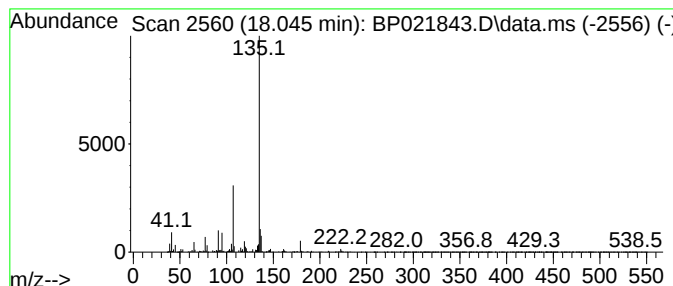
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Phenol, 4-(1,1,3,3-tetramet... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.045	2.47 ng/ul	457471	Phenanthrene-d10	17.210

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	78
2			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	78
3			Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	72
4			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	72
5			Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090424\
Data File : BP021843.D
Acq On : 05 Sep 2024 15:49
Operator : RC/JU
Sample : P3809-11
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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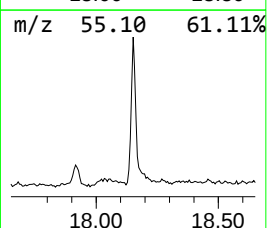
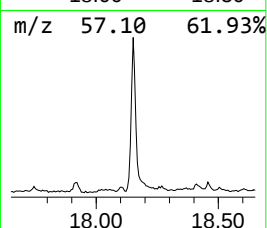
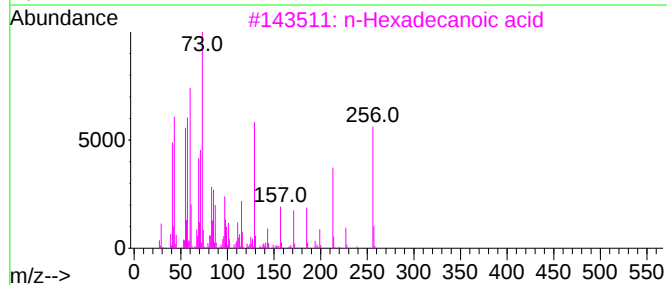
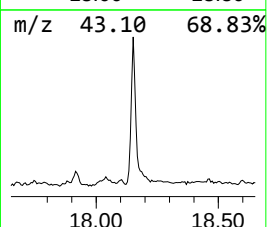
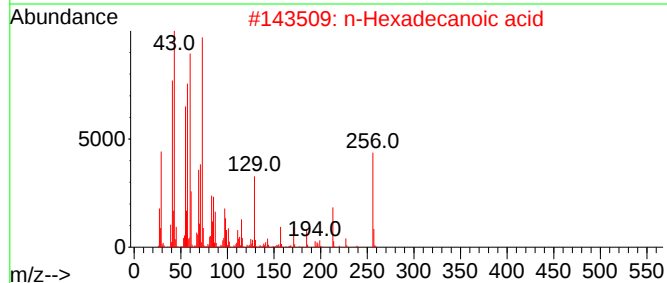
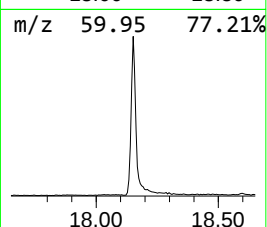
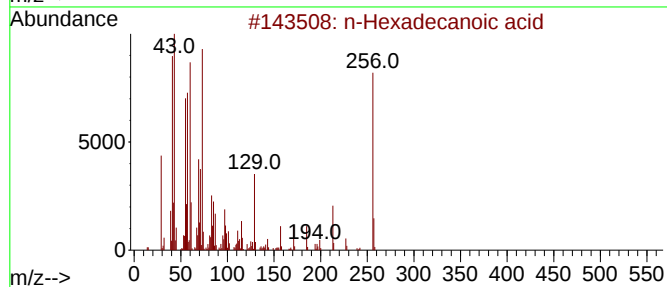
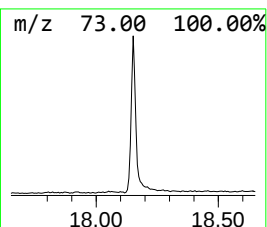
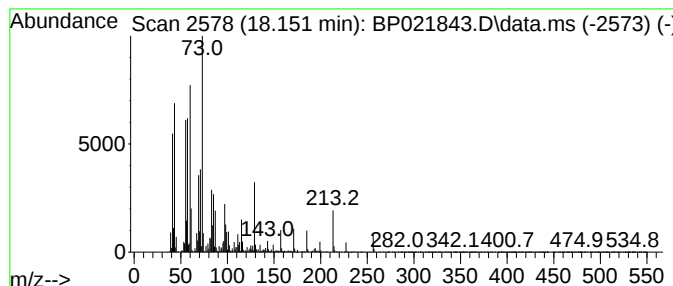
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.151	10.55 ng/ul	1955100	Phenanthrene-d10	17.210

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	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090424\
Data File : BP021843.D
Acq On : 05 Sep 2024 15:49
Operator : RC/JU
Sample : P3809-11
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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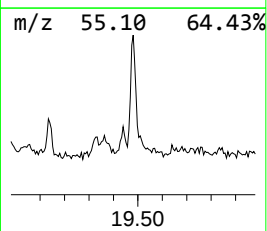
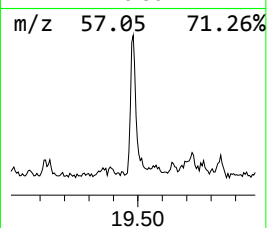
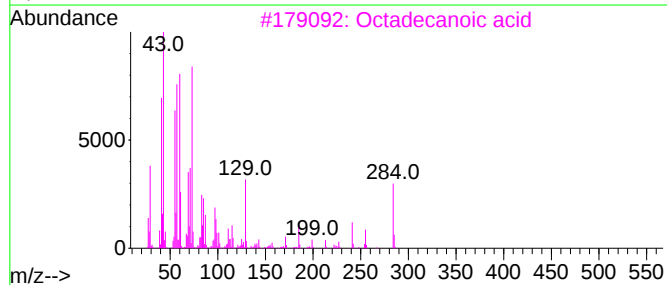
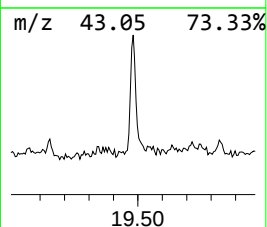
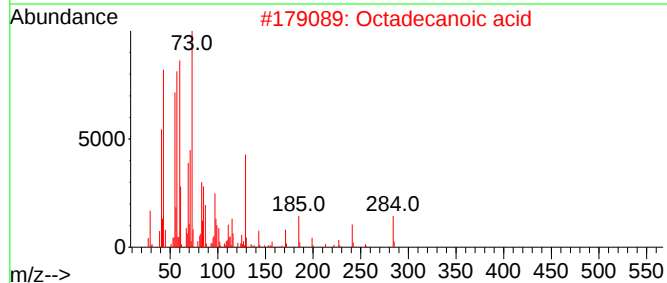
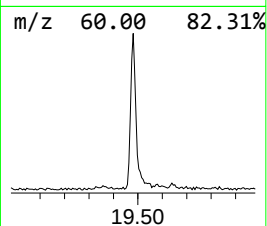
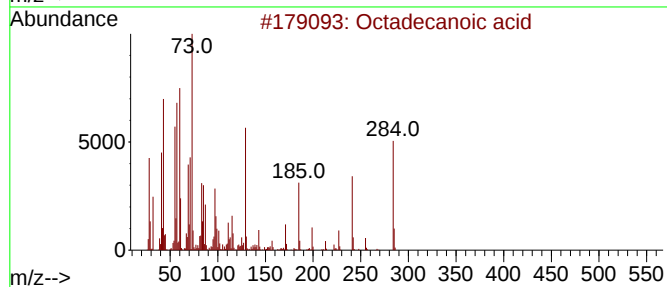
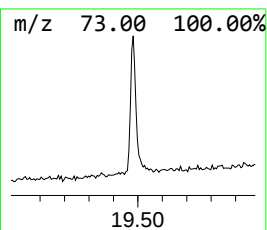
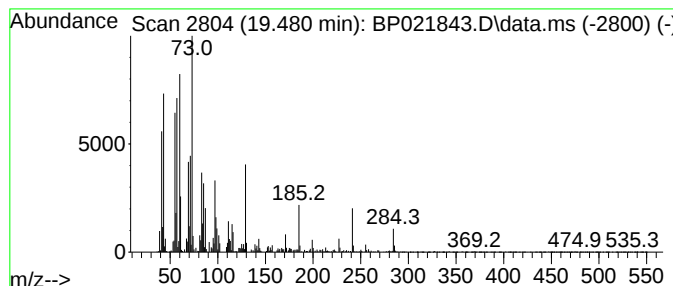
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.480	3.22 ng/ul	726105	Chrysene-d12	21.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	98
3			Octadecanoic acid	284	C18H36O2	000057-11-4	95
4			Octadecanoic acid	284	C18H36O2	000057-11-4	95
5			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090424\
Data File : BP021843.D
Acq On : 05 Sep 2024 15:49
Operator : RC/JU
Sample : P3809-11
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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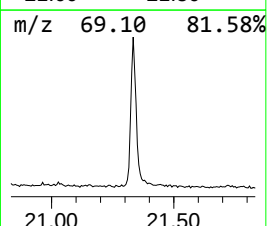
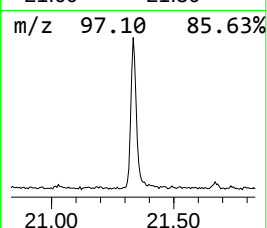
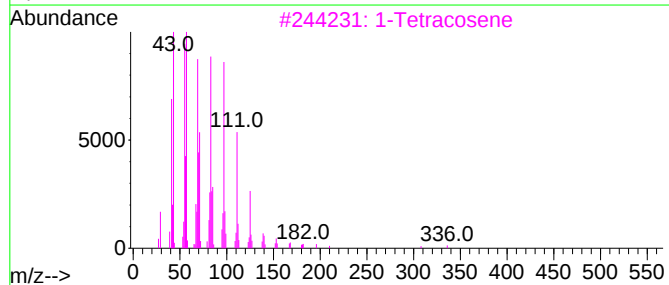
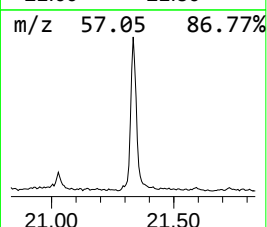
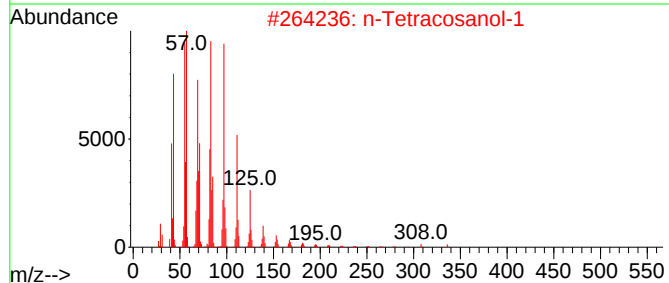
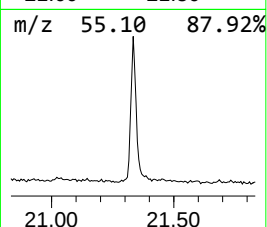
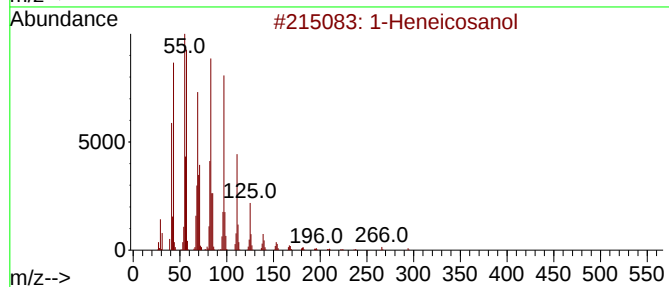
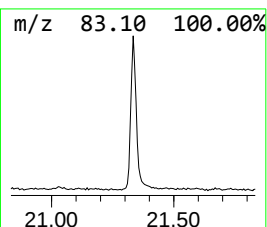
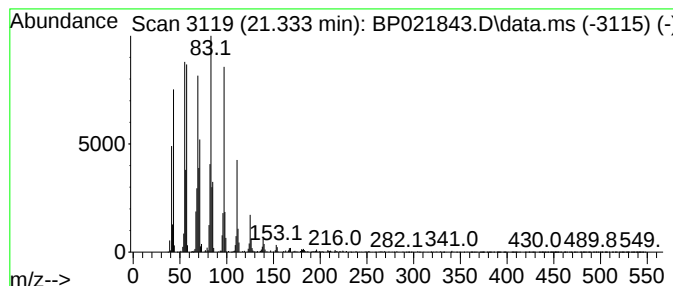
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 1-Heneicosanol Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.333	6.37 ng/ul	1434570	Chrysene-d12	21.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosanol			312	C21H44O	015594-90-8	94
2	n-Tetracosanol-1			354	C24H50O	000506-51-4	91
3	1-Tetracosene			336	C24H48	010192-32-2	91
4	Bromoacetic acid, hexadecyl ester			362	C18H35BrO2	005454-48-8	91
5	Bromoacetic acid, octadecyl ester			390	C20H39BrO2	018992-03-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090424\
Data File : BP021843.D
Acq On : 05 Sep 2024 15:49
Operator : RC/JU
Sample : P3809-11
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YE3E5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.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
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Ethanol, 2-(2-e...	7.369	2.2	ng/ul	158714	1	7.769	1435640	20.0
1-Phenoxypropan...	11.293	7.1	ng/ul	731983	2	10.552	2071780	20.0
unknown-01	14.334	2.7	ng/ul	366001	3	14.404	2737670	20.0
Pentadecanoic a...	17.916	2.3	ng/ul	432699	4	17.210	3705830	20.0
Phenol, 4-(1,1,...	18.045	2.5	ng/ul	457471	4	17.210	3705830	20.0
n-Hexadecanoic ...	18.151	10.6	ng/ul	1955100	4	17.210	3705830	20.0
Octadecanoic acid	19.480	3.2	ng/ul	726105	5	21.669	4506670	20.0
1-Heneicosanol	21.333	6.4	ng/ul	1434570	5	21.669	4506670	20.0