

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
 Data File : BP021681.D
 Acq On : 27 Aug 2024 21:56
 Operator : RC/JU
 Sample : P3675-18
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCXZ2

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: BP021681.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.017	3	5	23	rVB	3424088	4243998	63.77%	4.385%
2	3.193	31	35	42	rVB3	17654	26619	0.40%	0.028%
3	3.276	45	49	54	rVB	57364	70220	1.06%	0.073%
4	3.540	89	94	101	rBV	243194	331525	4.98%	0.343%
5	3.687	115	119	130	rBV	621488	862807	12.96%	0.891%
6	4.017	171	175	182	rVB	30793	44290	0.67%	0.046%
7	4.670	282	286	292	rBV	17621	27673	0.42%	0.029%
8	4.923	325	329	334	rVB	13837	21824	0.33%	0.023%
9	5.164	361	370	374	rBV5	11620	23677	0.36%	0.024%
10	5.893	489	494	501	rVB	24574	37165	0.56%	0.038%
11	6.381	573	577	584	rVB8	7403	14968	0.22%	0.015%
12	6.752	632	640	644	rBV	39092	78805	1.18%	0.081%
13	6.787	644	646	653	rVV2	26827	50940	0.77%	0.053%
14	6.958	667	675	686	rBV	899156	1440675	21.65%	1.488%
15	7.122	696	703	712	rVB	776223	1241720	18.66%	1.283%
16	7.328	731	738	744	rBV	884326	1409070	21.17%	1.456%
17	7.393	744	749	763	rVB	165856	277931	4.18%	0.287%
18	7.787	809	816	823	rBV	1382761	2332896	35.05%	2.410%
19	7.852	823	827	836	rVB3	22138	39156	0.59%	0.040%
20	8.487	927	935	944	rBV	780719	1267290	19.04%	1.309%
21	8.940	1004	1012	1021	rBV	846457	1462732	21.98%	1.511%
22	9.105	1035	1040	1048	rVB	10972	19249	0.29%	0.020%
23	9.499	1102	1107	1113	rVB	97035	145428	2.19%	0.150%
24	9.658	1122	1134	1146	rBV	666293	1123927	16.89%	1.161%
25	9.811	1152	1160	1169	rBV2	34012	93038	1.40%	0.096%
26	9.893	1169	1174	1183	rVV2	10328	28939	0.43%	0.030%
27	10.199	1215	1226	1244	rBV	958632	1927108	28.95%	1.991%
28	10.352	1248	1252	1260	rVB9	11372	25111	0.38%	0.026%
29	10.575	1281	1290	1302	rBV	1960641	3330476	50.04%	3.441%
30	10.705	1304	1312	1330	rVV	768886	1409306	21.17%	1.456%
31	11.116	1376	1382	1392	rVB3	36175	72201	1.08%	0.075%
32	11.322	1410	1417	1432	rBV	208563	510267	7.67%	0.527%
33	12.069	1539	1544	1554	rBV3	14912	32019	0.48%	0.033%
34	12.169	1554	1561	1565	rBV3	23027	45235	0.68%	0.047%
35	12.581	1626	1631	1635	rBV6	8834	17552	0.26%	0.018%
36	12.799	1663	1668	1673	rBV7	11733	18930	0.28%	0.020%

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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_P\Methods\SFAM-EPA-BP082324.MA.M
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37	13.210	1733	1738	1743	rVV9	12677	26694	0.40%	0.028%
38	13.328	1751	1758	1764	rVV7	27713	61378	0.92%	0.063%
39	13.387	1764	1768	1774	rVV7	15968	25155	0.38%	0.026%
40	13.563	1793	1798	1802	rVV2	12807	24494	0.37%	0.025%
41	13.763	1826	1832	1837	rBV2	52452	83812	1.26%	0.087%
42	13.834	1837	1844	1859	rVV	1820224	2840162	42.67%	2.934%
43	14.122	1882	1893	1911	rBV	2122982	3462125	52.02%	3.577%
44	14.287	1913	1921	1925	rVV7	19738	43772	0.66%	0.045%
45	14.334	1925	1929	1935	rVV9	14541	28601	0.43%	0.030%
46	14.434	1937	1946	1955	rVV	2884545	4542844	68.26%	4.693%
47	14.510	1956	1959	1965	rVB5	19798	31815	0.48%	0.033%
48	14.646	1973	1982	1983	rBV4	2872834	4595115	69.04%	4.747%
49	14.663	1983	1985	2000	rVB	3480164	4470822	67.17%	4.619%
50	14.857	2014	2018	2023	rVB4	32692	42942	0.65%	0.044%
51	14.951	2030	2034	2047	rVB10	23126	62929	0.95%	0.065%
52	15.193	2069	2075	2080	rBV2	26748	49685	0.75%	0.051%
53	15.251	2082	2085	2090	rVB7	11433	17896	0.27%	0.018%
54	15.357	2098	2103	2108	rVV2	34099	66847	1.00%	0.069%
55	15.440	2108	2117	2129	rVV	2907425	4434615	66.63%	4.582%
56	15.551	2129	2136	2145	rVV	616295	965945	14.51%	0.998%
57	15.634	2145	2150	2155	rVV7	32636	72034	1.08%	0.074%
58	15.687	2156	2159	2166	rVB6	21223	37537	0.56%	0.039%
59	15.893	2190	2194	2198	rBV7	21795	34010	0.51%	0.035%
60	16.016	2211	2215	2216	rBV4	15137	19678	0.30%	0.020%
61	16.110	2227	2231	2235	rVB5	15775	25684	0.39%	0.027%
62	16.187	2241	2244	2249	rVB6	16357	20379	0.31%	0.021%
63	16.245	2251	2254	2257	rVB5	12814	14392	0.22%	0.015%
64	16.357	2269	2273	2278	rVV8	15794	33725	0.51%	0.035%
65	16.434	2280	2286	2291	rVV3	38013	74129	1.11%	0.077%
66	16.622	2312	2318	2323	rVB4	42401	68708	1.03%	0.071%
67	16.928	2367	2370	2377	rVB8	26462	44650	0.67%	0.046%
68	17.163	2406	2410	2414	rVV4	82193	134060	2.01%	0.139%
69	17.222	2414	2420	2427	rVV	3541015	5539857	83.24%	5.723%
70	17.328	2427	2438	2449	rVV2	3080951	4783882	71.88%	4.942%
71	17.416	2449	2453	2460	rVB9	25446	44925	0.67%	0.046%
72	17.528	2469	2472	2475	rBV4	15737	20426	0.31%	0.021%
73	17.616	2482	2487	2498	rBV4	45415	113056	1.70%	0.117%
74	17.934	2537	2541	2549	rVB5	45958	79509	1.19%	0.082%
75	18.016	2550	2555	2564	rBV	134647	244161	3.67%	0.252%
76	18.104	2566	2570	2575	rVB8	37024	63612	0.96%	0.066%

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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 >
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77	18.169	2575	2581	2596	rBV	1767970	2981807	44.80%	3.081%
78	18.504	2636	2638	2651	rVB8	54616	138122	2.08%	0.143%
79	18.657	2660	2664	2669	rVB3	83324	108357	1.63%	0.112%
80	18.851	2694	2697	2703	rBV2	41393	54532	0.82%	0.056%
81	19.057	2729	2732	2738	rBV3	58701	90061	1.35%	0.093%
82	19.151	2744	2748	2752	rVB3	93026	100425	1.51%	0.104%
83	19.210	2755	2758	2761	rBV4	34187	41292	0.62%	0.043%
84	19.245	2761	2764	2769	rVB3	57687	67054	1.01%	0.069%
85	19.339	2769	2780	2781	rBV3	97299	229688	3.45%	0.237%
86	19.381	2781	2787	2792	rVB	229667	406050	6.10%	0.420%
87	19.422	2792	2794	2799	rVB2	66032	73529	1.10%	0.076%
88	19.498	2802	2807	2819	rBV	1043936	1687172	25.35%	1.743%
89	19.692	2834	2840	2847	rBV2	3828900	5819379	87.44%	6.012%
90	19.975	2884	2888	2894	rBV3	148674	192134	2.89%	0.199%
91	20.169	2919	2921	2926	rVB	87480	101623	1.53%	0.105%
92	20.257	2933	2936	2940	rBV3	174205	262446	3.94%	0.271%
93	20.633	2996	3000	3004	rBV	303208	319725	4.80%	0.330%
94	21.133	3081	3085	3090	rBV	297994	408071	6.13%	0.422%
95	21.351	3117	3122	3133	rBV	727260	1241763	18.66%	1.283%
96	21.675	3170	3177	3184	rBV2	3782573	6562961	98.61%	6.780%
97	22.633	3334	3340	3353	rVB	658916	1529613	22.98%	1.580%
98	24.257	3612	3616	3622	rBV	239699	461961	6.94%	0.477%
99	24.857	3708	3718	3728	rBV2	2089689	5807593	87.26%	6.000%
100	25.092	3749	3758	3772	rVB3	2269387	6655593	100.00%	6.876%

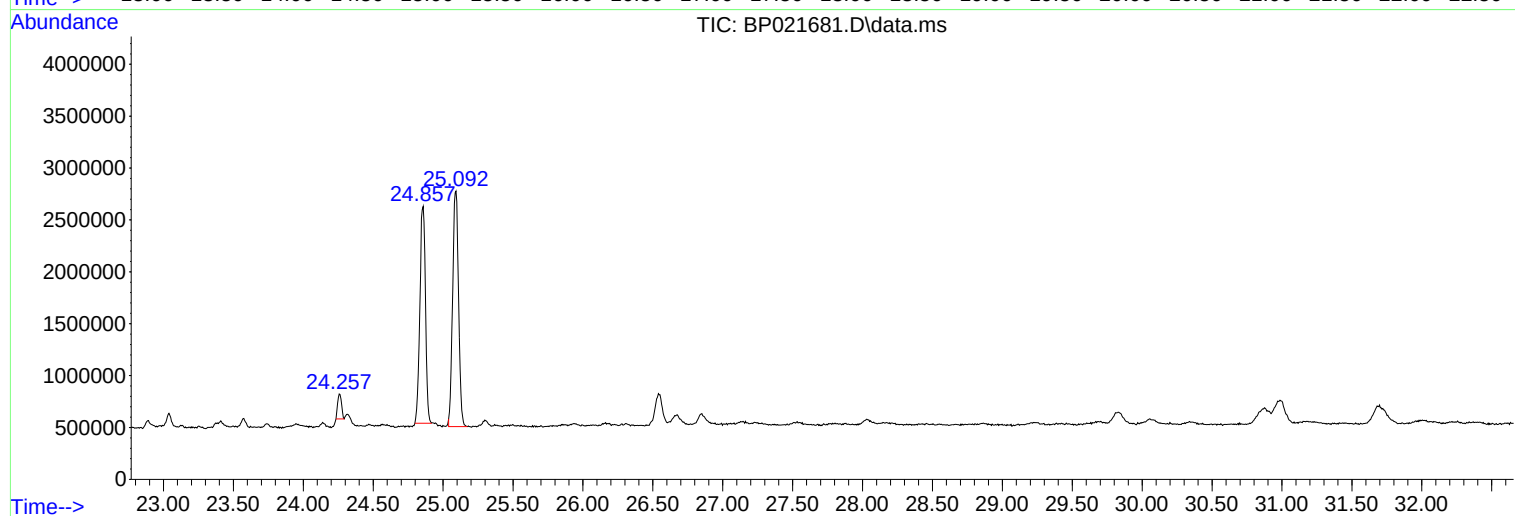
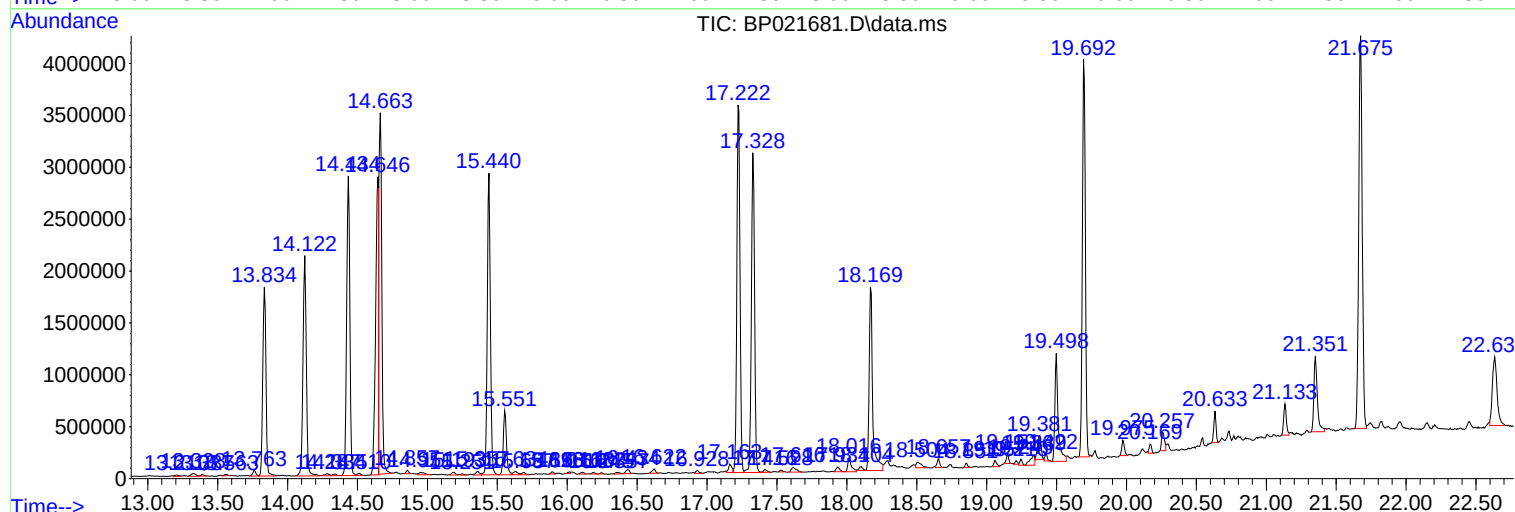
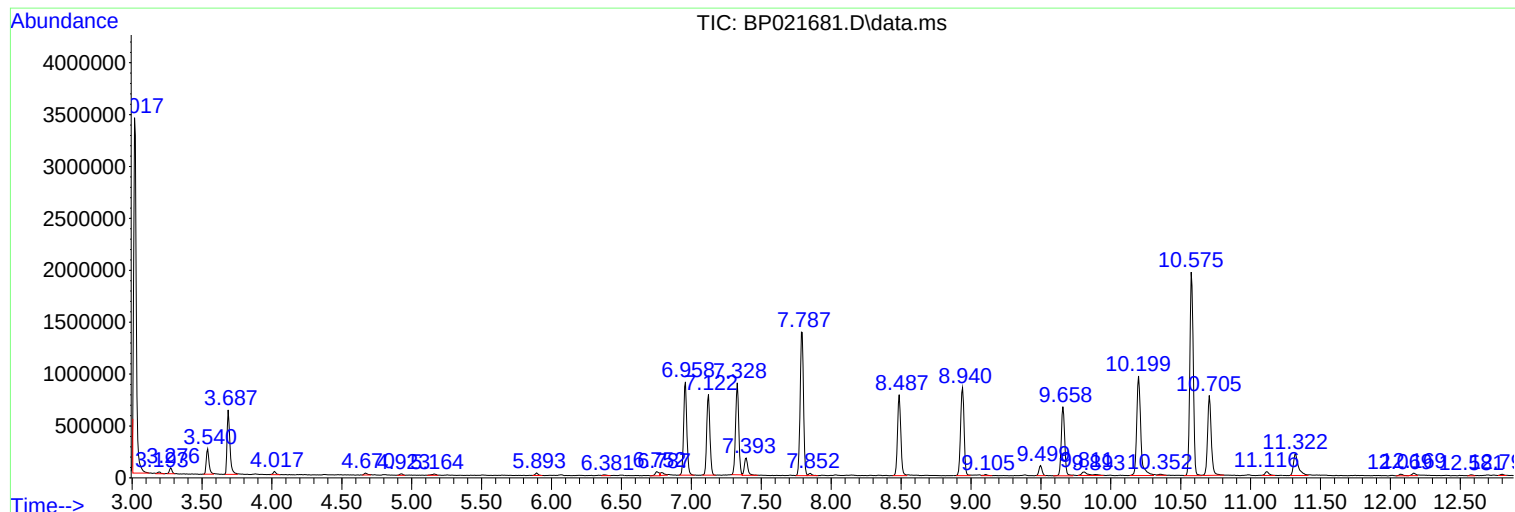
Sum of corrected areas: 96791780

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Instrument :
BNA_P
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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



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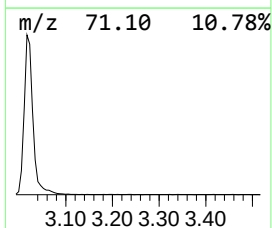
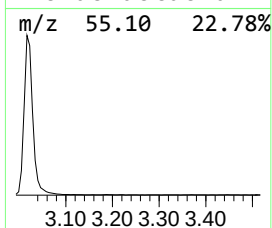
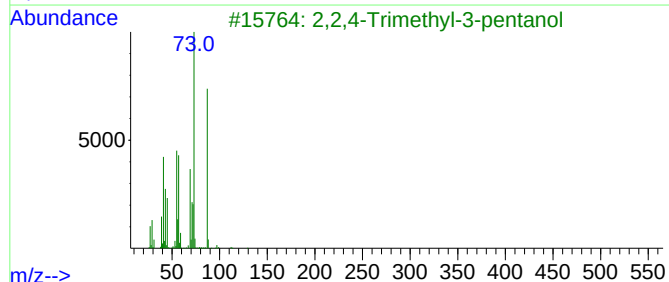
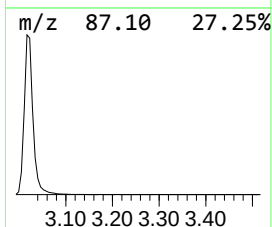
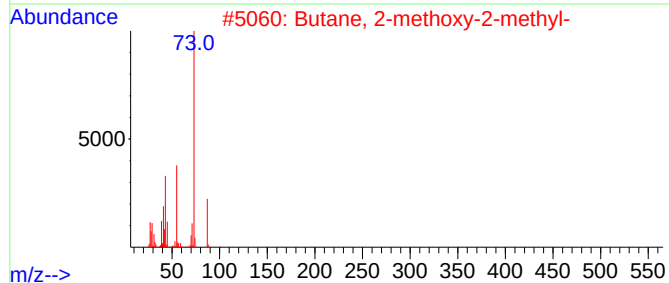
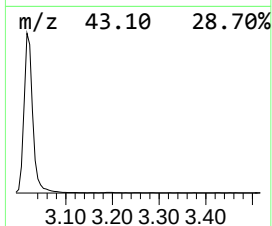
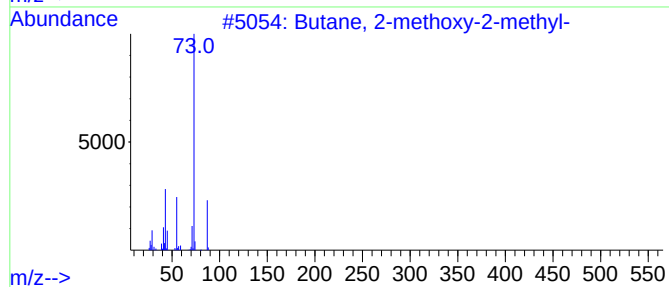
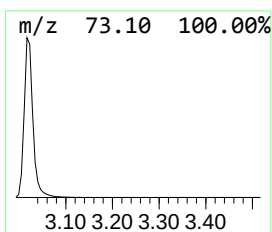
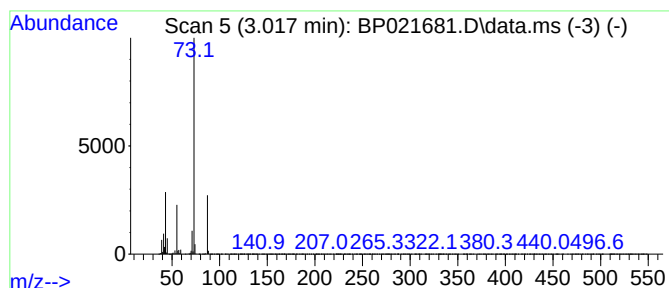
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.017	36.38 ng/ul	4244000	1,4-Dichlorobenzene-d4	7.793

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83		
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78		
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40		
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17		
5	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9		



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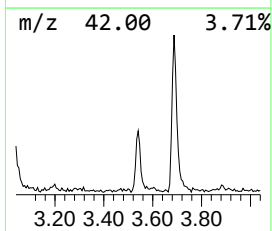
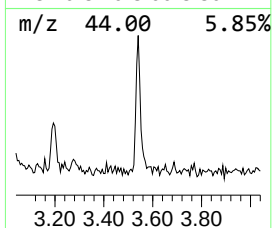
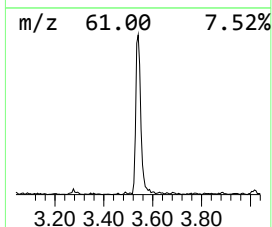
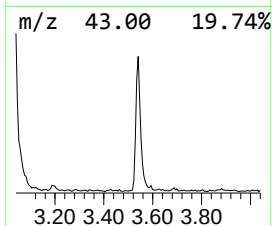
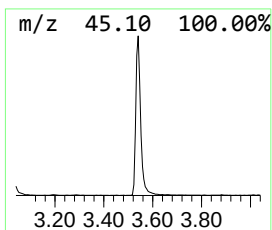
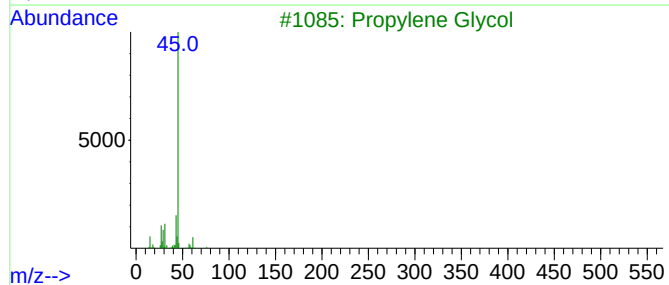
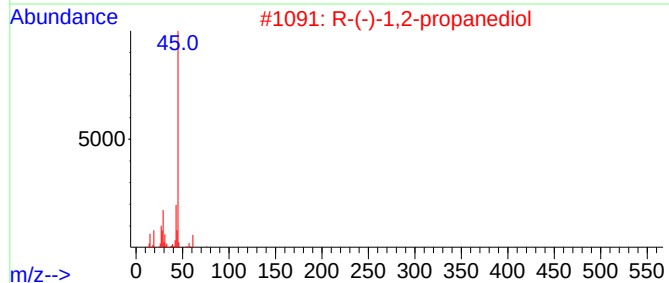
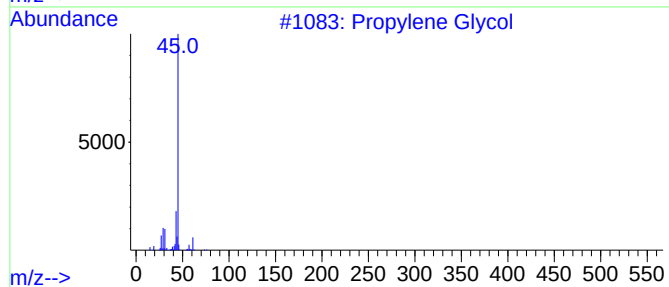
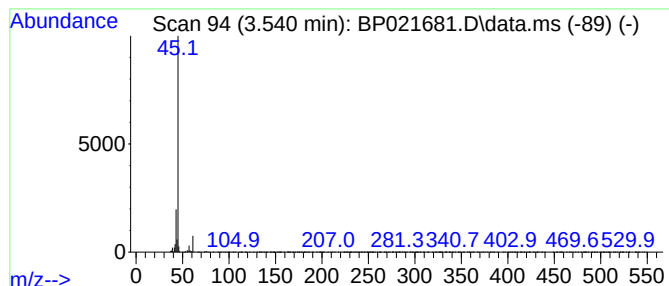
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Propylene Glycol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.540	2.84 ng/ul	331525	1,4-Dichlorobenzene-d4	7.793

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	90
2			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	78
3			Propylene Glycol	76	C3H8O2	000057-55-6	64
4			Isopropyl Alcohol	60	C3H8O	000067-63-0	56
5			Propanoic acid, 2-hydroxy-, meth...	104	C4H8O3	000547-64-8	40



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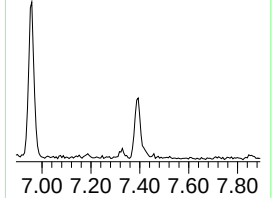
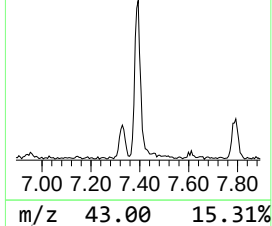
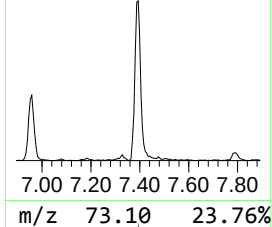
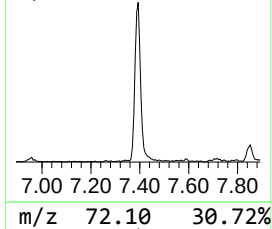
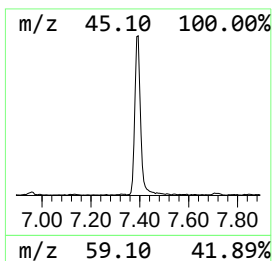
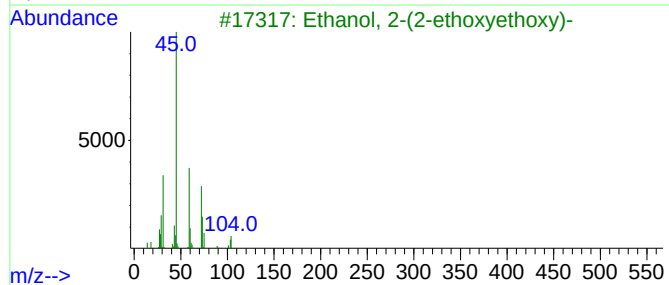
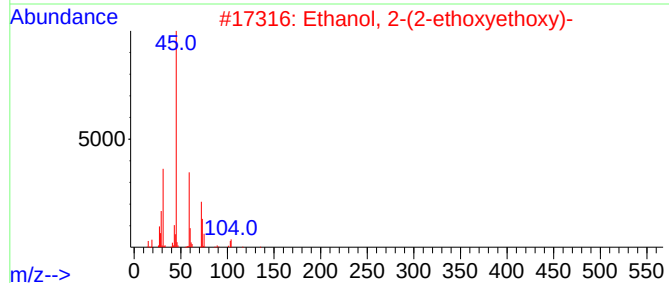
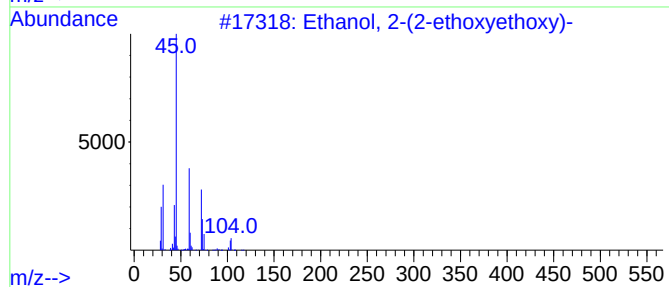
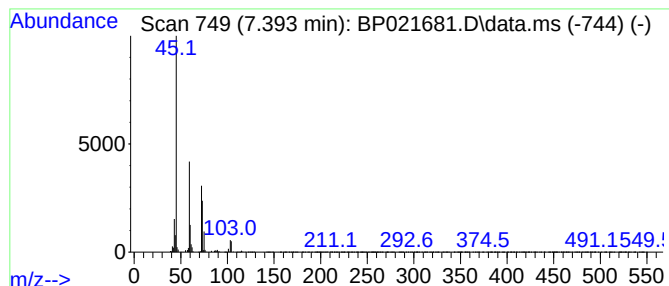
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Ethanol, 2-(2-ethoxyethoxy)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.393	2.38 ng/ul	277931	1,4-Dichlorobenzene-d4	7.793

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	78
3			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	72
4			Ethanol, 2-(2-ethoxyethoxy)-	134	C6H14O3	000111-90-0	72
5			Diethyl carbitol	162	C8H18O3	000112-36-7	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021681.D
Acq On : 27 Aug 2024 21:56
Operator : RC/JU
Sample : P3675-18
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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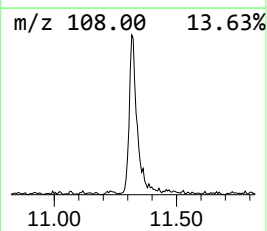
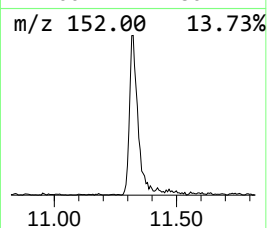
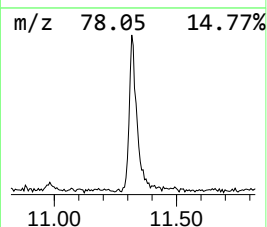
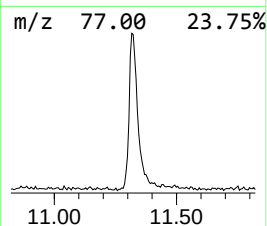
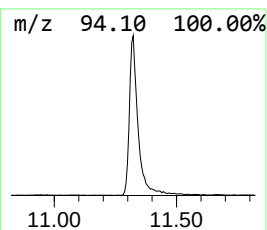
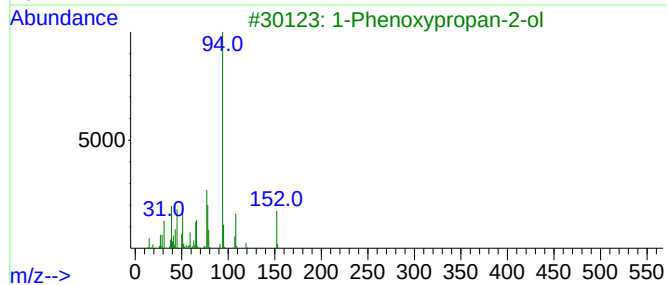
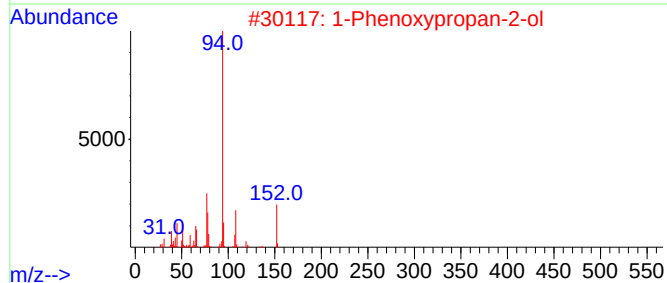
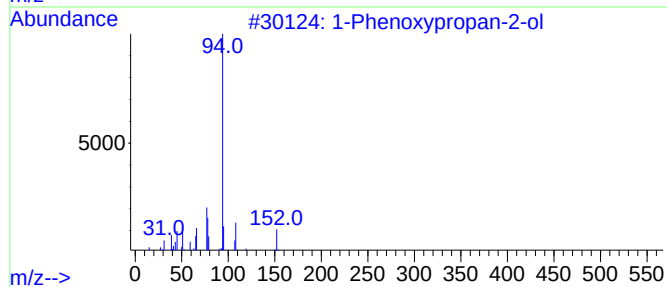
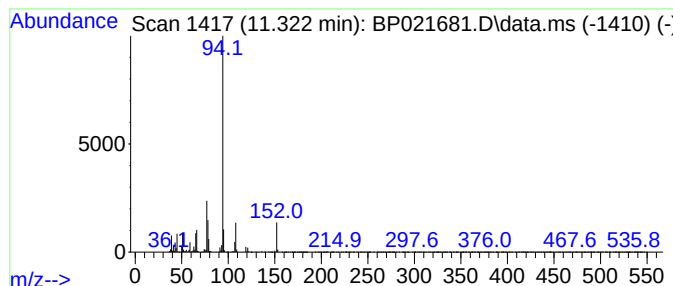
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Phenoxypropan-2-ol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.322	3.06 ng/ul	510267	Naphthalene-d8	10.575

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	94
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
4			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	87
5			1-Propanol, 2-phenoxy-	152	C9H12O2	004169-04-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021681.D
Acq On : 27 Aug 2024 21:56
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Sample : P3675-18
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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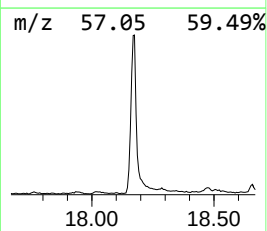
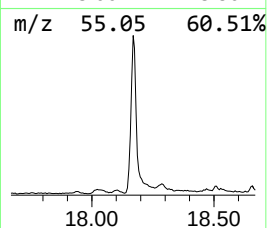
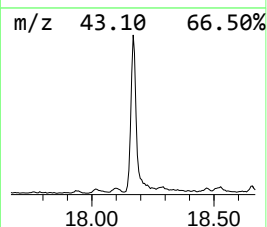
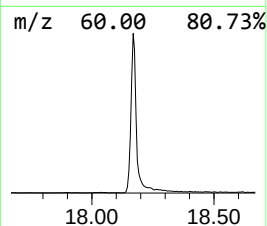
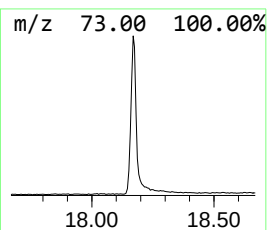
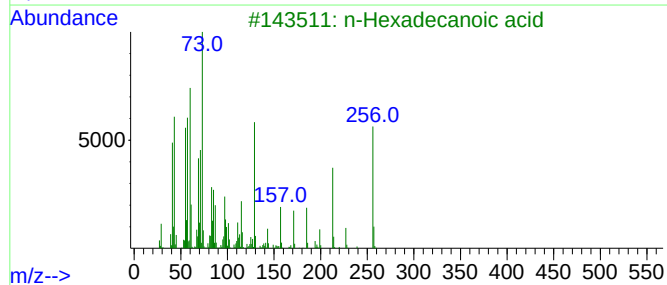
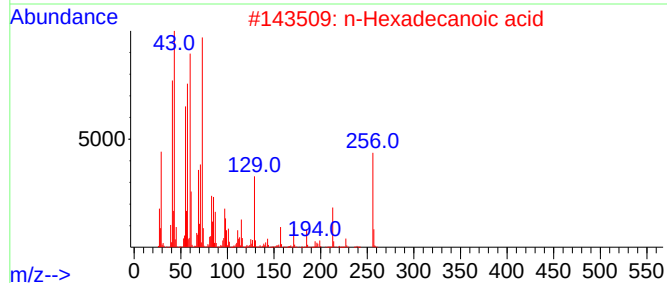
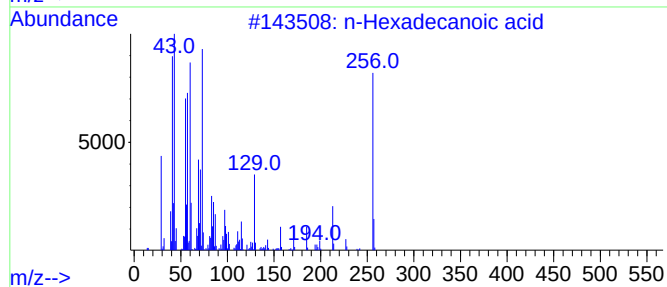
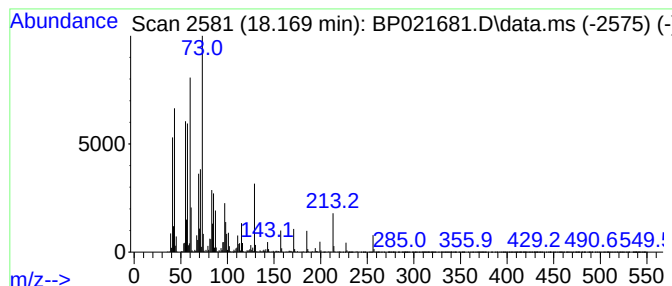
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.169	10.76 ng/ul	2981810	Phenanthrene-d10	17.222

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\BP082624\
Data File : BP021681.D
Acq On : 27 Aug 2024 21:56
Operator : RC/JU
Sample : P3675-18
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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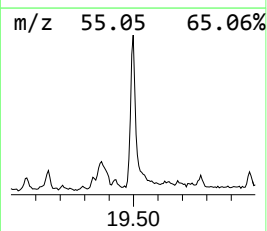
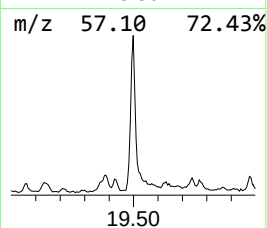
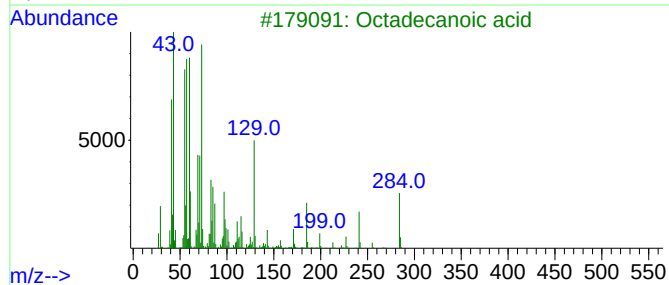
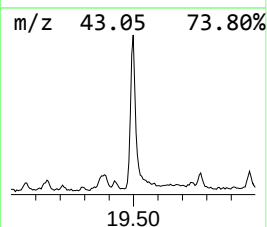
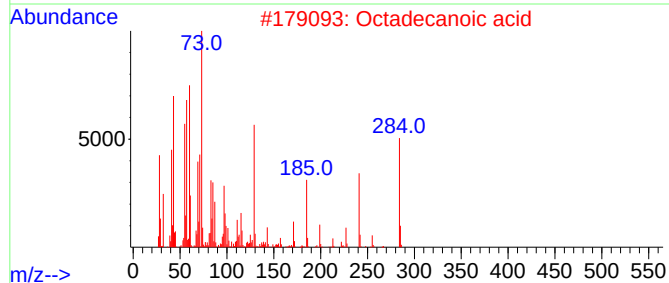
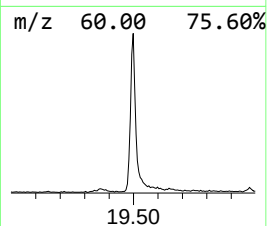
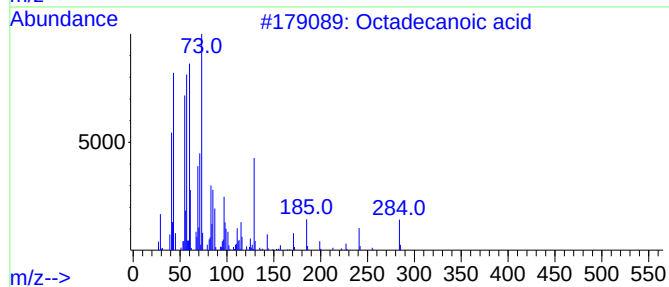
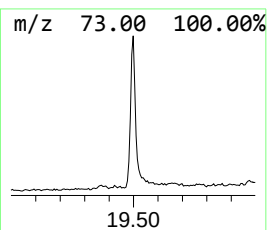
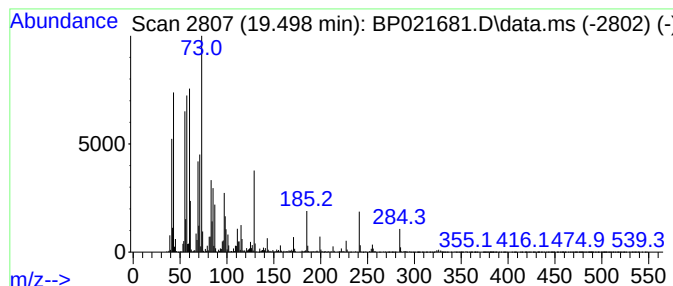
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.498	5.14 ng/ul	1687170	Chrysene-d12	21.675

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	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	97
4			Octadecanoic acid	284	C18H36O2	000057-11-4	96
5			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021681.D
Acq On : 27 Aug 2024 21:56
Operator : RC/JU
Sample : P3675-18
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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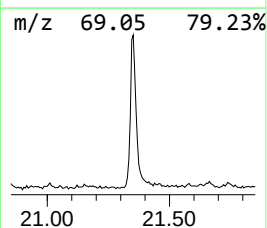
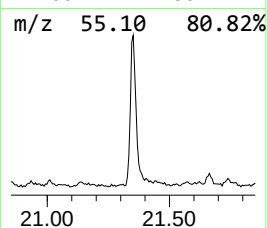
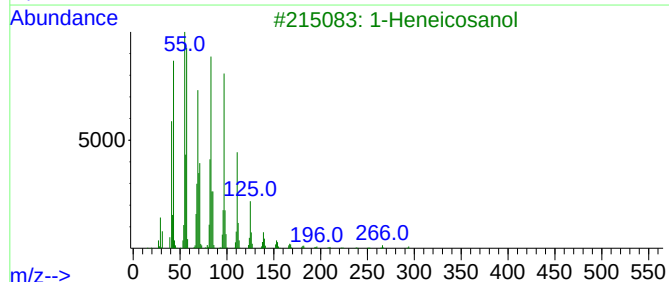
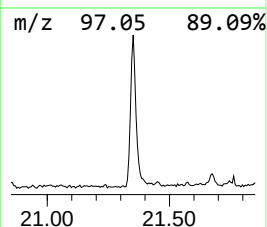
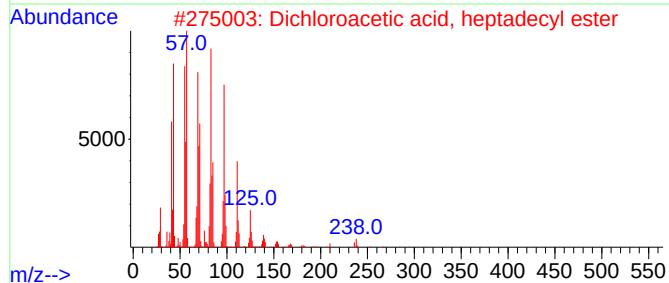
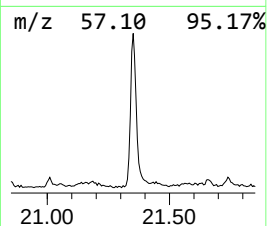
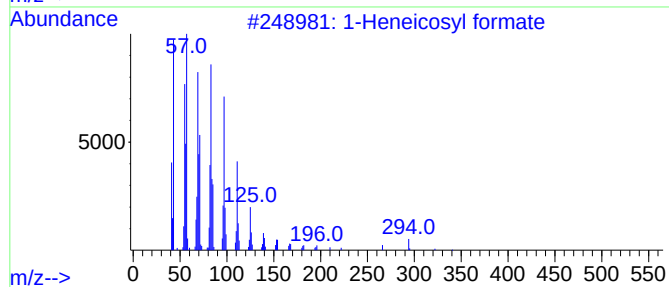
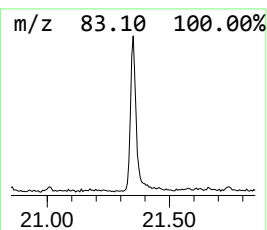
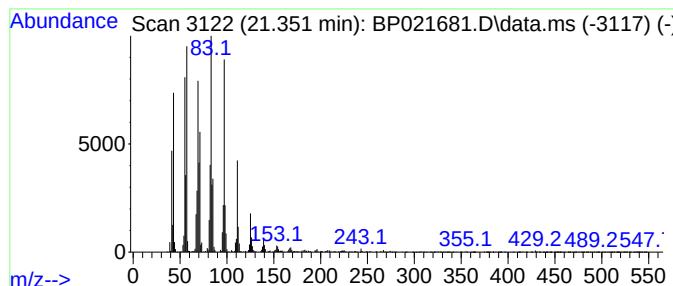
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1-Heneicosyl formate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.351	3.78 ng/ul	1241760	Chrysene-d12	21.675

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosyl formate	340	C22H44O2	077899-03-7	94
2			Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	91
3			1-Heneicosanol	312	C21H44O	015594-90-8	91
4			n-Nonadecanol-1	284	C19H40O	001454-84-8	91
5			Behenic alcohol	326	C22H46O	000661-19-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021681.D
Acq On : 27 Aug 2024 21:56
Operator : RC/JU
Sample : P3675-18
Misc :
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ClientSampleId :
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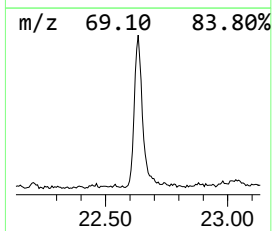
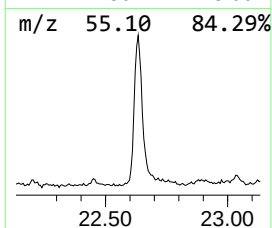
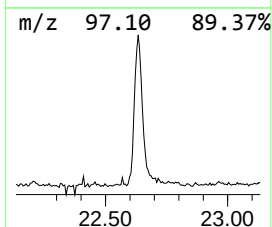
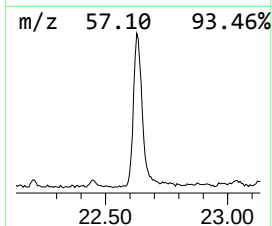
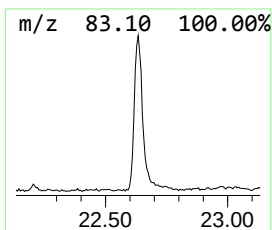
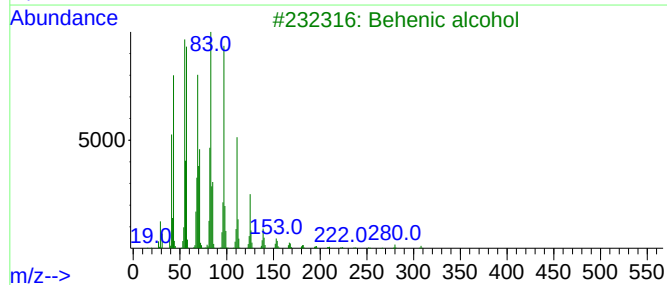
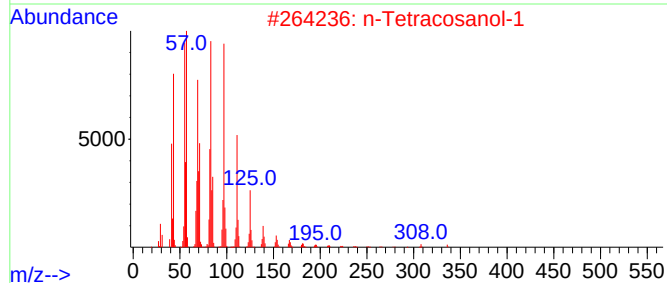
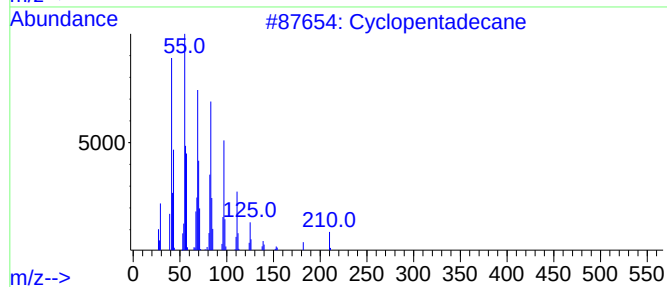
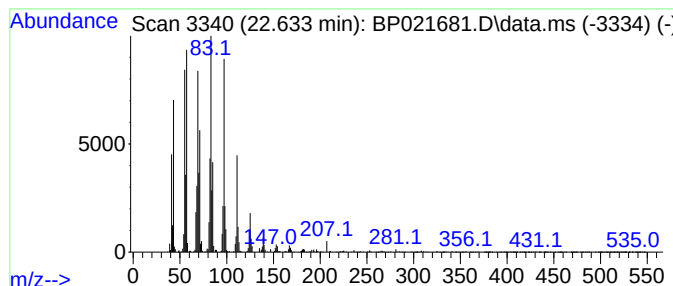
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Cyclic22.633 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.633	4.66 ng/ul	1529610	Chrysene-d12	21.675

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentadecane	210	C15H30	000295-48-7	93
2			n-Tetracosanol-1	354	C24H50O	000506-51-4	91
3			Behenic alcohol	326	C22H46O	000661-19-8	91
4			1-Heptacosanol	396	C27H56O	002004-39-9	91
5			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
Data File : BP021681.D
Acq On : 27 Aug 2024 21:56
Operator : RC/JU
Sample : P3675-18
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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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\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Propylene Glycol	3.540	2.8	ng/ul	331525	1	7.793	2332900	20.0
Ethanol, 2-(2-e...	7.393	2.4	ng/ul	277931	1	7.793	2332900	20.0
1-Phenoxypropan...	11.322	3.1	ng/ul	510267	2	10.575	3330480	20.0
n-Hexadecanoic ...	18.169	10.8	ng/ul	2981810	4	17.222	5539860	20.0
Octadecanoic acid	19.498	5.1	ng/ul	1687170	5	21.675	6562960	20.0
1-Heneicosyl fo...	21.351	3.8	ng/ul	1241760	5	21.675	6562960	20.0
(DEL) Alkane: C...	22.633	4.7	ng/ul	1529610	5	21.675	6562960	20.0