

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042224\
 Data File : BP020077.D
 Acq On : 22 Apr 2024 15:12
 Operator : MA/JU
 Sample : P2104-14
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AC5

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-BP041924.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP020077.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.228	21	24	30	rVB2	18580	28489	1.26%	0.109%
2	3.499	67	70	87	rVB	41302	76516	3.39%	0.292%
3	3.634	89	93	115	rVB	202207	348794	15.47%	1.331%
4	3.828	124	126	128	rBV3	2643	2287	0.10%	0.009%
5	3.928	140	143	151	rVB4	6440	9787	0.43%	0.037%
6	5.675	435	440	441	rBV5	1797	2375	0.11%	0.009%
7	6.693	610	613	620	rBV9	2982	7348	0.33%	0.028%
8	6.828	629	636	653	rBV	302261	566627	25.12%	2.162%
9	6.993	658	664	674	rVV	227385	384931	17.07%	1.469%
10	7.093	679	681	685	rVV4	2324	2445	0.11%	0.009%
11	7.181	689	696	712	rVV	291244	525267	23.29%	2.004%
12	7.334	716	722	726	rVV9	3240	7798	0.35%	0.030%
13	7.640	767	774	785	rBV	353990	607661	26.94%	2.318%
14	8.258	874	879	881	rBV6	1965	2853	0.13%	0.011%
15	8.346	887	894	909	rBV	342074	686738	30.45%	2.620%
16	8.799	964	971	985	rBV	282161	533476	23.65%	2.035%
17	9.152	1026	1031	1038	rVB2	6902	13055	0.58%	0.050%
18	9.375	1064	1069	1074	rBV7	6066	10721	0.48%	0.041%
19	9.510	1084	1092	1104	rBV	259442	475466	21.08%	1.814%
20	9.957	1163	1168	1172	rBV7	3943	7111	0.32%	0.027%
21	10.052	1176	1184	1202	rBV	338667	708559	31.42%	2.703%
22	10.252	1213	1218	1219	rBV5	3488	4508	0.20%	0.017%
23	10.416	1239	1246	1257	rBV	451769	816997	36.23%	3.117%
24	10.575	1264	1273	1293	rBV	333680	702765	31.16%	2.681%
25	11.199	1372	1379	1394	rBV2	17046	53380	2.37%	0.204%
26	11.846	1483	1489	1495	rVB2	2268	4074	0.18%	0.016%
27	12.040	1513	1522	1531	rVB6	4980	11327	0.50%	0.043%
28	12.904	1664	1669	1672	rBV6	2583	4936	0.22%	0.019%
29	13.122	1702	1706	1712	rVB8	3618	6647	0.29%	0.025%
30	13.198	1714	1719	1725	rVV4	9544	18150	0.80%	0.069%
31	13.281	1727	1733	1739	rVB4	2562	4071	0.18%	0.016%
32	13.740	1803	1811	1823	rBV	717913	1091048	48.38%	4.163%
33	14.010	1847	1857	1869	rBV	802038	1300600	57.67%	4.962%
34	14.228	1892	1894	1900	rVB7	2146	3034	0.13%	0.012%
35	14.328	1903	1911	1921	rBV	717168	1113806	49.39%	4.249%
36	14.545	1939	1948	1949	rBV	45325	58659	2.60%	0.224%

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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-BP041924.MA.M
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37	14.592	1949	1956	1973	rVB2	169761	501770	22.25%	1.914%
38	14.775	1983	1987	1994	rVB8	11324	22666	1.01%	0.086%
39	15.151	2049	2051	2055	rVB4	2542	2785	0.12%	0.011%
40	15.216	2058	2062	2067	rVB7	2833	4850	0.22%	0.019%
41	15.357	2079	2086	2102	rBV	1063041	1638490	72.65%	6.251%
42	15.504	2105	2111	2124	rVV	172895	315707	14.00%	1.205%
43	15.604	2124	2128	2137	rVB	7273	15347	0.68%	0.059%
44	15.792	2158	2160	2164	rVB5	1805	2529	0.11%	0.010%
45	15.945	2179	2186	2192	rBV7	8235	16003	0.71%	0.061%
46	16.110	2211	2214	2217	rBV4	5504	6668	0.30%	0.025%
47	16.298	2240	2246	2250	rBV8	3715	8343	0.37%	0.032%
48	16.575	2289	2293	2298	rBV8	5316	8800	0.39%	0.034%
49	16.710	2310	2316	2319	rBV6	5196	9647	0.43%	0.037%
50	16.939	2350	2355	2357	rBV4	7262	9090	0.40%	0.035%
51	17.116	2380	2385	2388	rBV6	8557	11238	0.50%	0.043%
52	17.175	2388	2395	2406	rVB	805130	1394025	61.81%	5.319%
53	17.286	2407	2414	2424	rVV	1172583	1829803	81.13%	6.981%
54	17.404	2427	2434	2440	rBV	6104	17246	0.76%	0.066%
55	17.510	2447	2452	2454	rBV6	9034	15738	0.70%	0.060%
56	17.598	2465	2467	2475	rVB9	9108	14964	0.66%	0.057%
57	17.722	2486	2488	2493	rVB5	5381	5887	0.26%	0.022%
58	17.904	2514	2519	2523	rBV	28178	39074	1.73%	0.149%
59	18.004	2533	2536	2539	rBV5	3565	4934	0.22%	0.019%
60	18.145	2554	2560	2570	rBV	309825	481900	21.37%	1.839%
61	18.451	2608	2612	2614	rBV5	12188	17692	0.78%	0.067%
62	19.045	2709	2713	2719	rBV9	14055	23128	1.03%	0.088%
63	19.228	2741	2744	2749	rBV4	16003	24710	1.10%	0.094%
64	19.328	2758	2761	2763	rBV2	17717	17676	0.78%	0.067%
65	19.492	2783	2789	2797	rBV2	115001	225697	10.01%	0.861%
66	19.680	2813	2821	2832	rVB	1546726	2255277	100.00%	8.604%
67	20.292	2922	2925	2928	rVB	37679	41282	1.83%	0.158%
68	20.692	2986	2993	3009	rBV	1331553	1764316	78.23%	6.731%
69	21.357	3102	3106	3114	rVB	252871	384369	17.04%	1.466%
70	21.680	3155	3161	3172	rBV	939906	1431896	63.49%	5.463%
71	24.851	3692	3700	3714	rVB2	720311	2038096	90.37%	7.776%
72	25.080	3731	3739	3750	rVB	525254	1406596	62.37%	5.367%

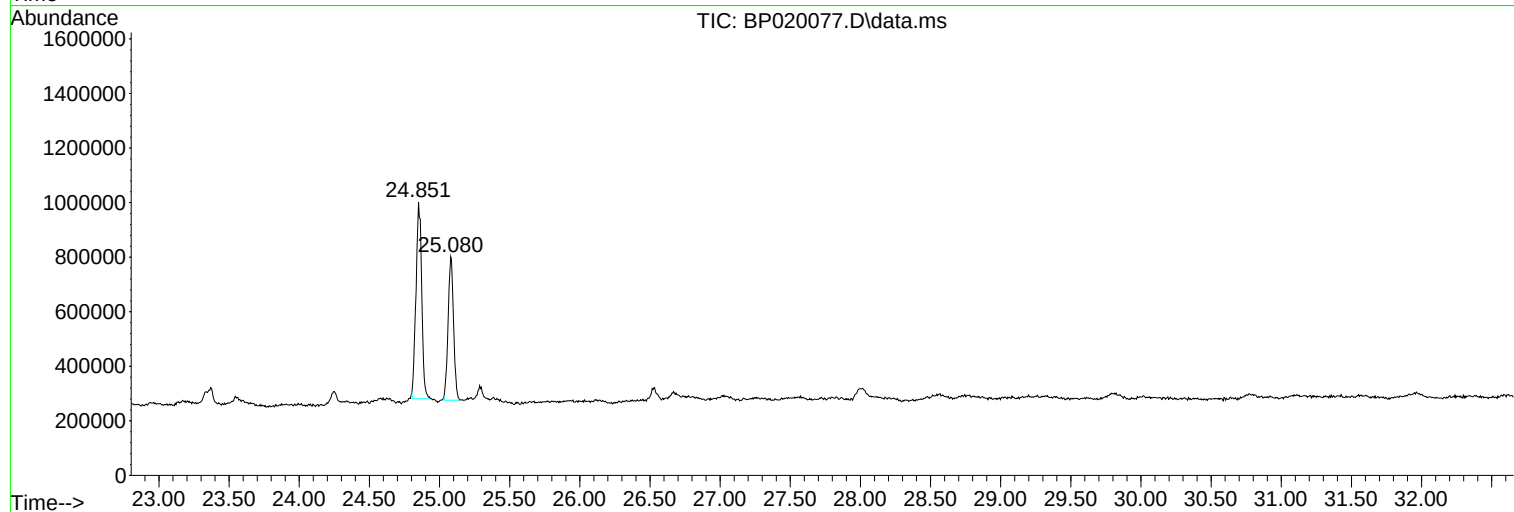
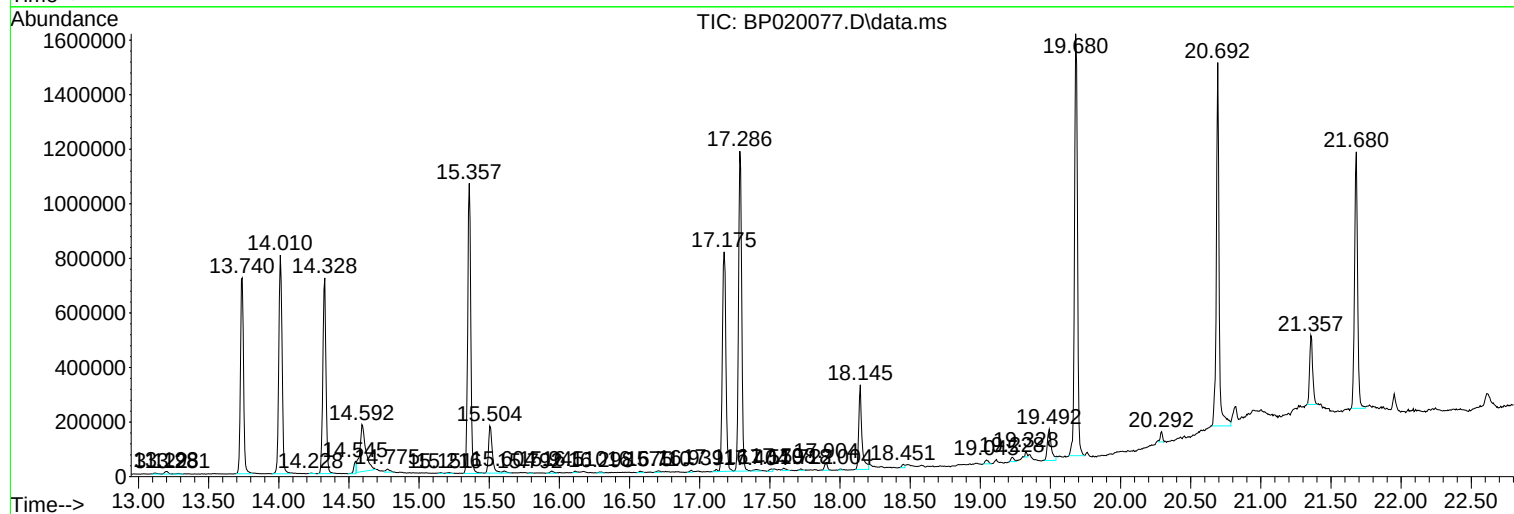
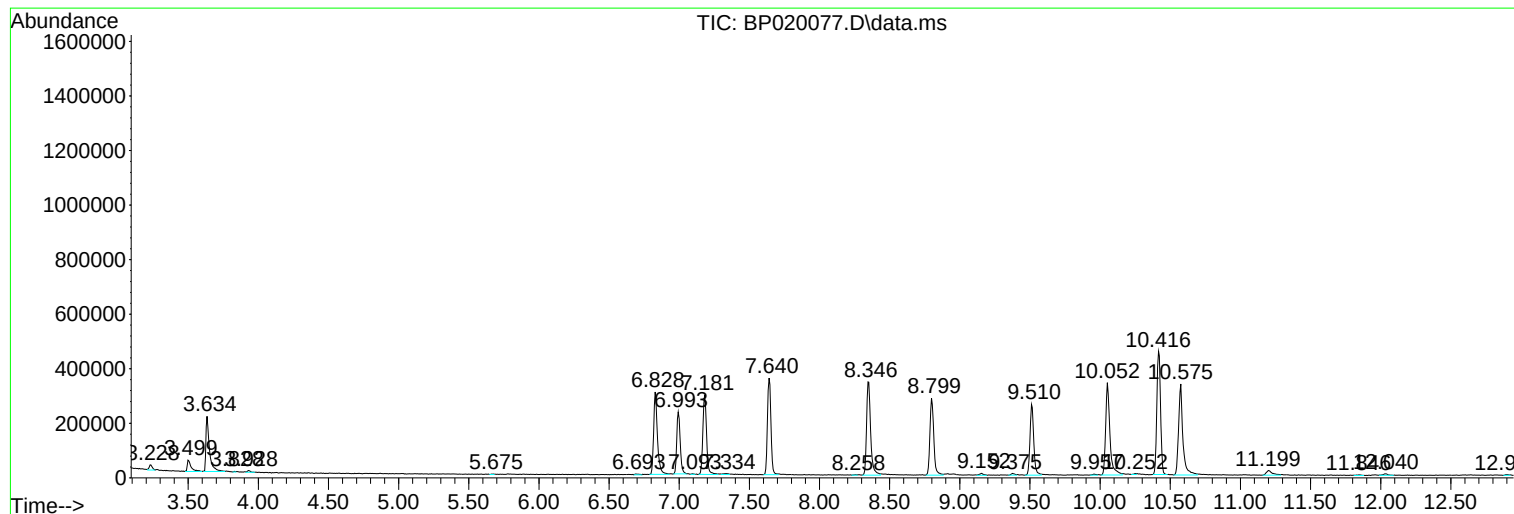
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041924.MA.M
Quant Title : SVOA CALIBRATION

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



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Data File : BP020077.D
Acq On : 22 Apr 2024 15:12
Operator : MA/JU
Sample : P2104-14
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AC5

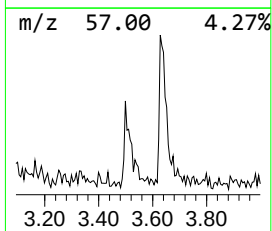
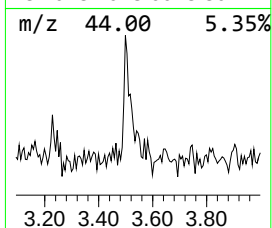
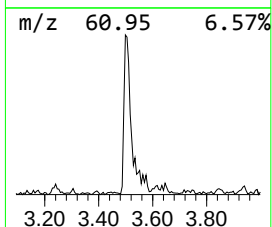
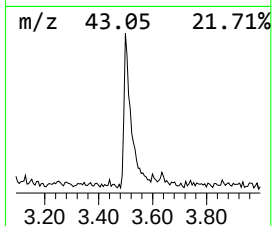
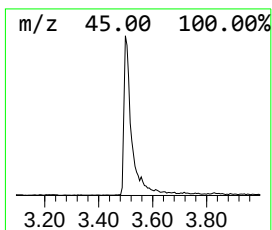
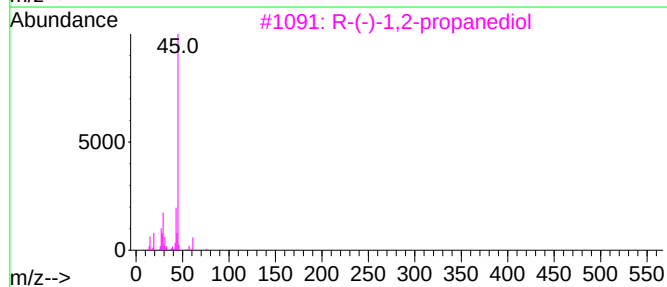
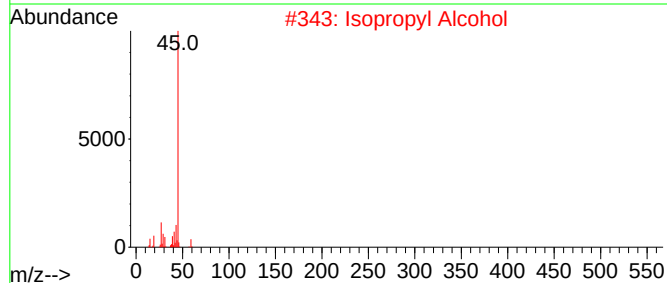
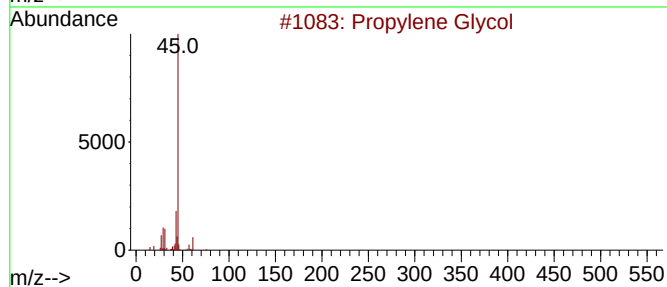
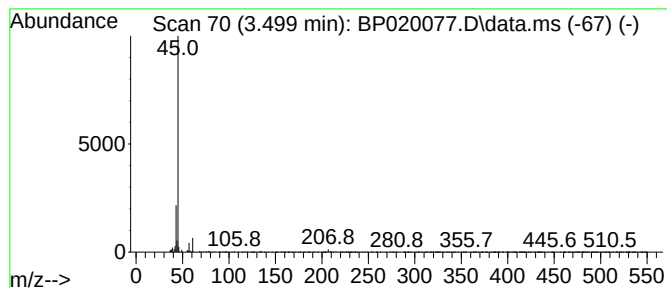
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 Propylene Glycol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.499	2.52 ng/ul	76516	1,4-Dichlorobenzene-d4	7.640

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propylene Glycol	76	C3H8O2	000057-55-6	72
2			Isopropyl Alcohol	60	C3H8O	000067-63-0	50
3			R-(-)-1,2-propanediol	76	C3H8O2	004254-14-2	9
4			(S)-Isopropyl lactate	132	C6H12O3	063697-00-7	9
5			Propylene Glycol	76	C3H8O2	000057-55-6	9



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ALS Vial : 6 Sample Multiplier: 1

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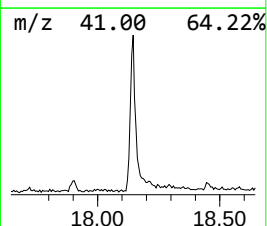
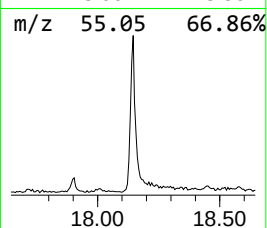
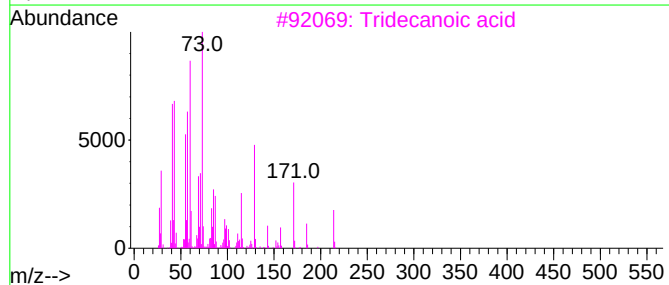
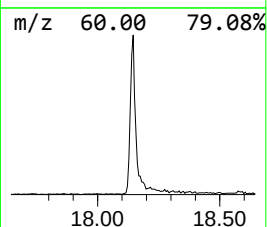
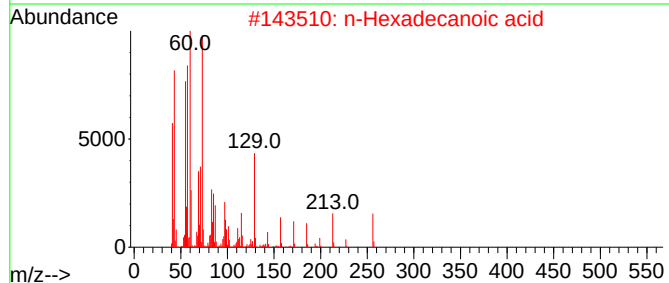
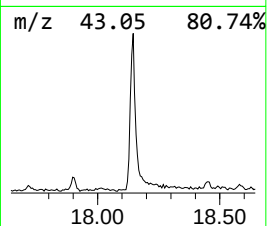
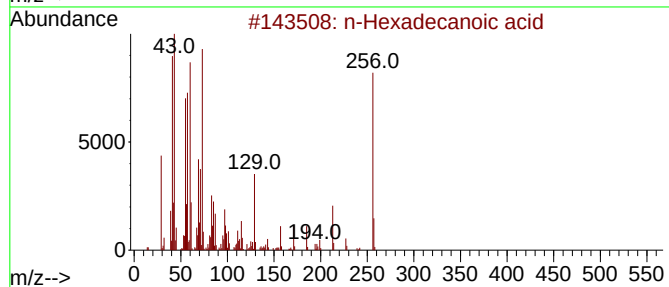
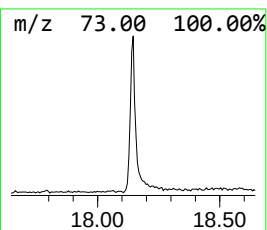
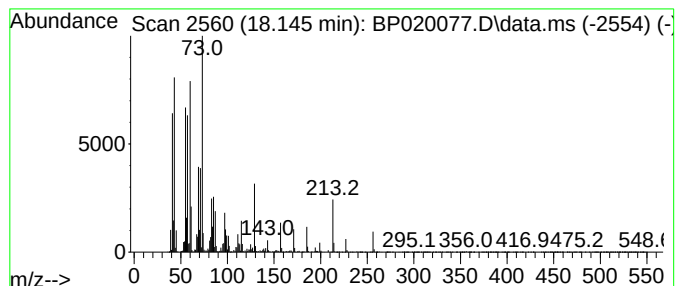
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.145	6.91 ng/ul	481900	Phenanthrene-d10	17.175

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	97
3			Tridecanoic acid	214	C13H26O2	000638-53-9	96
4			Tridecanoic acid	214	C13H26O2	000638-53-9	93
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	91



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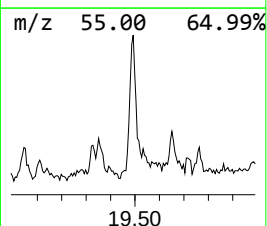
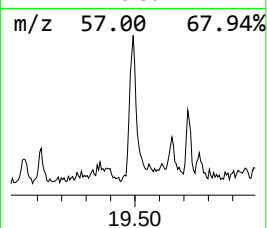
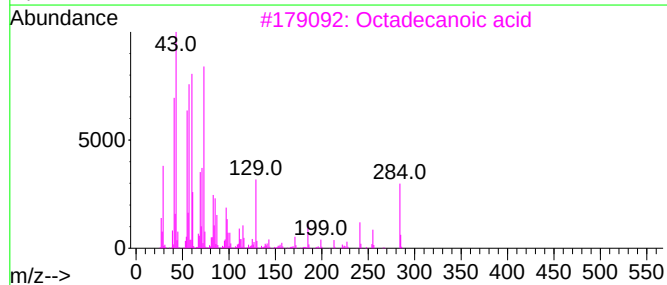
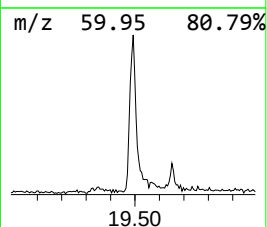
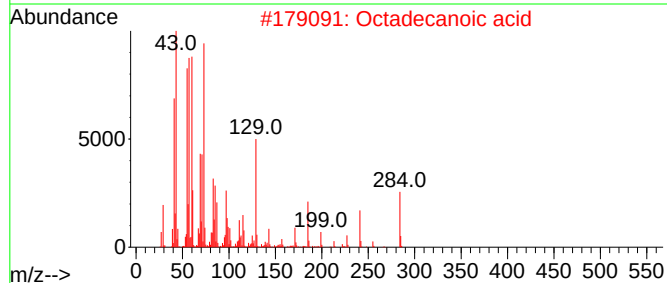
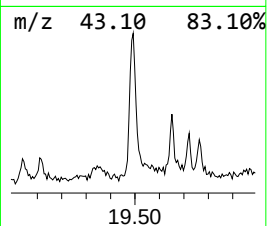
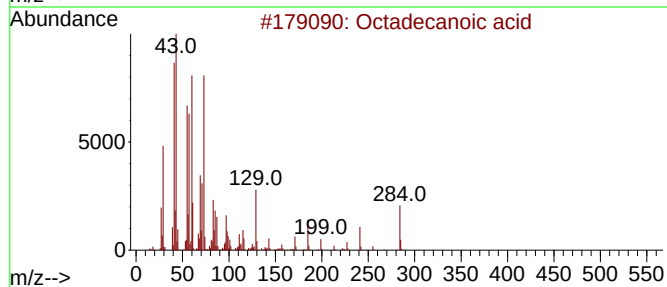
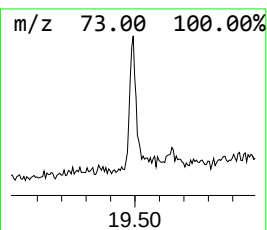
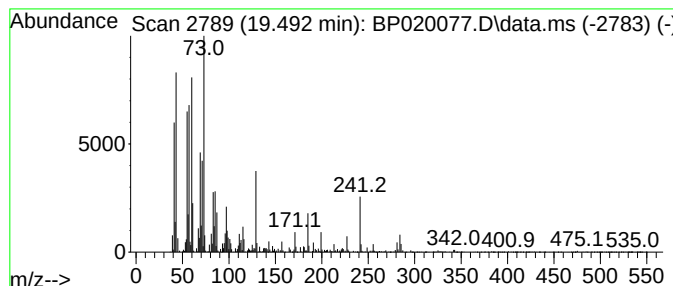
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.492	3.15 ng/ul	225697	Chrysene-d12	21.680

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	98
2			Octadecanoic acid	284	C18H36O2	000057-11-4	92
3			Octadecanoic acid	284	C18H36O2	000057-11-4	91
4			Pentadecanoic acid	242	C15H30O2	001002-84-2	90
5			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	80



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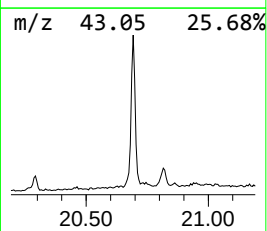
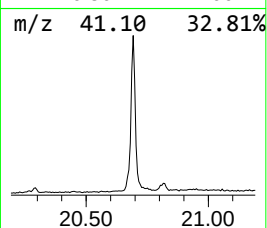
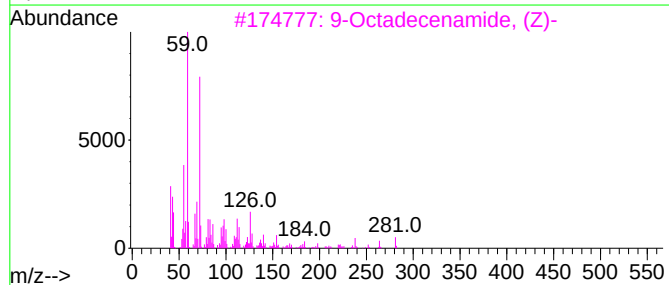
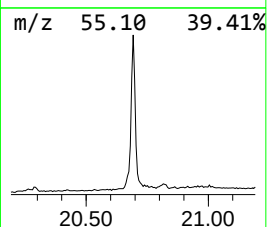
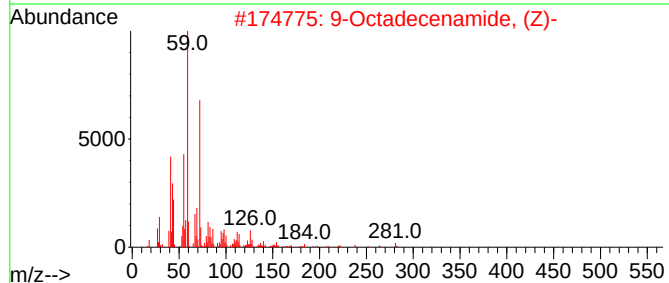
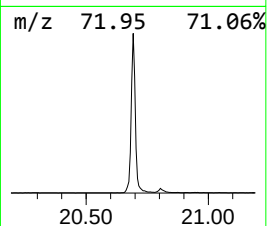
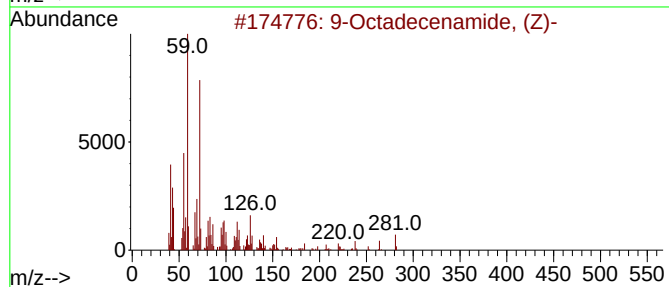
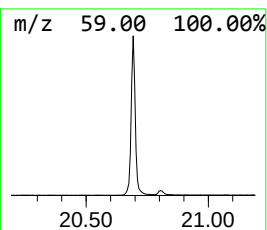
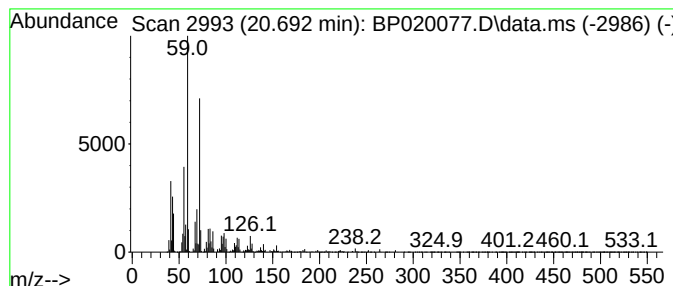
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041924.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 9-Octadecenamide, (Z)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.692	24.64 ng/ul	1764320	Chrysene-d12	21.680

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	93
2			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	91
3			9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	91
4			Hexadecanamide	255	C16H33NO	000629-54-9	78
5			Octanamide	143	C8H17NO	000629-01-6	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042224\
Data File : BP020077.D
Acq On : 22 Apr 2024 15:12
Operator : MA/JU
Sample : P2104-14
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AC5

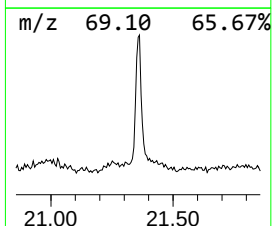
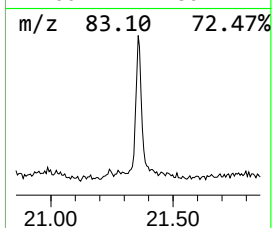
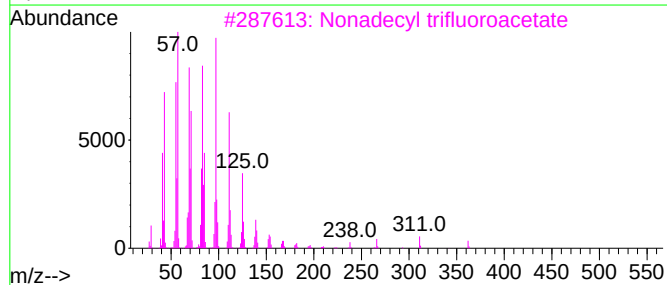
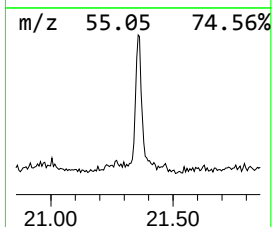
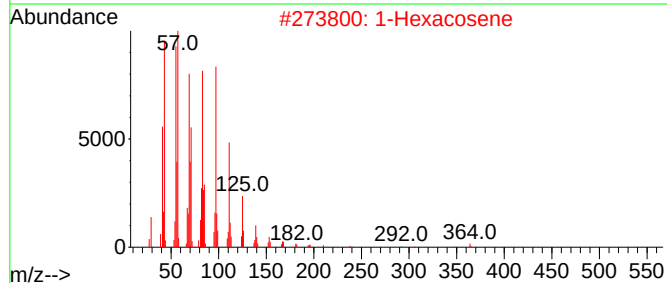
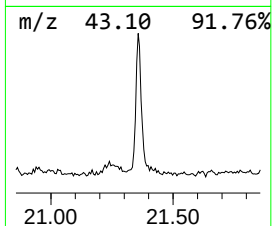
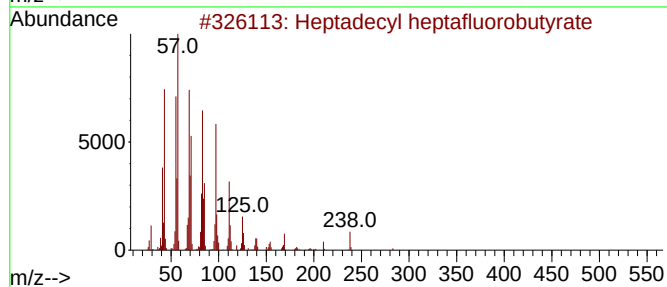
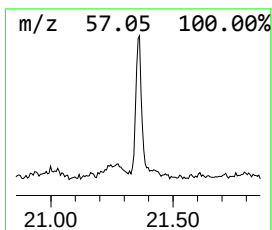
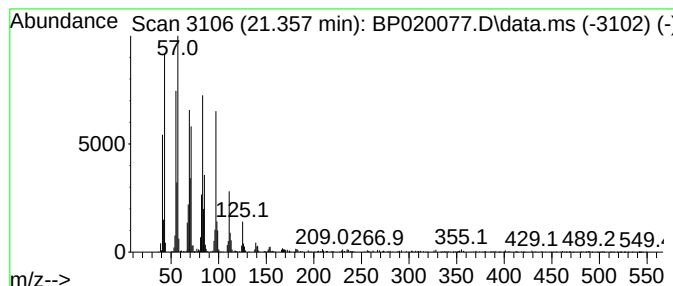
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Heptadecyl heptafluorobutyrate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.357	5.37 ng/ul	384369	Chrysene-d12	21.680

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
2			1-Hexacosene	364	C26H52	018835-33-1	94
3			Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	91
4			Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	91
5			17-Pentatriacontene	491	C35H70	006971-40-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042224\
Data File : BP020077.D
Acq On : 22 Apr 2024 15:12
Operator : MA/JU
Sample : P2104-14
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AC5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041924.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
Propylene Glycol	3.499	2.5	ng/ul	76516	1	7.640	607661	20.0
n-Hexadecanoic ...	18.145	6.9	ng/ul	481900	4	17.175	1394030	20.0
Octadecanoic acid	19.492	3.1	ng/ul	225697	5	21.680	1431900	20.0
9-Octadecenamid...	20.692	24.6	ng/ul	1764320	5	21.680	1431900	20.0
Heptadecyl hept...	21.357	5.4	ng/ul	384369	5	21.680	1431900	20.0