

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
 Data File : BP014225.D
 Acq On : 22 Mar 2023 19:54
 Operator : CG/JU
 Sample : 01977-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EYG88

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-BP030723.M
 Title : SVOA CALIBRATION

Signal : TIC: BP014225.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.405	33	37	47	rVB	52101	77974	0.93%	0.096%
2	3.699	83	87	96	rBV	34257	60728	0.73%	0.075%
3	3.840	107	111	131	rBV	586322	1020483	12.21%	1.253%
4	4.170	162	167	172	rVB2	9466	14067	0.17%	0.017%
5	5.364	366	370	378	rBV3	5810	10923	0.13%	0.013%
6	7.199	675	682	696	rBV	896412	1539690	18.42%	1.891%
7	7.364	704	710	723	rVB	748583	1181241	14.13%	1.450%
8	7.569	737	745	761	rBV	935764	1495865	17.90%	1.837%
9	8.040	818	825	833	rBV	514406	811927	9.71%	0.997%
10	8.758	939	947	967	rBV	1156398	1960689	23.46%	2.407%
11	9.216	1018	1025	1040	rBV	963336	1618267	19.36%	1.987%
12	9.369	1040	1051	1059	rVV	8763	23384	0.28%	0.029%
13	9.599	1085	1090	1095	rVB3	10148	16171	0.19%	0.020%
14	9.946	1140	1149	1162	rBV	834175	1453808	17.39%	1.785%
15	10.222	1185	1196	1197	rBV10	4979	10525	0.13%	0.013%
16	10.375	1216	1222	1232	rVB10	6491	14487	0.17%	0.018%
17	10.487	1232	1241	1262	rBV	1158992	2133878	25.53%	2.620%
18	10.863	1298	1305	1316	rBV	765790	1285122	15.38%	1.578%
19	11.005	1318	1329	1358	rVV	885904	1798310	21.52%	2.208%
20	12.451	1569	1575	1585	rVB2	20816	37222	0.45%	0.046%
21	13.293	1711	1718	1721	rBV7	4839	8852	0.11%	0.011%
22	13.493	1747	1752	1756	rBV5	14725	24733	0.30%	0.030%
23	13.569	1761	1765	1771	rVB2	17273	27991	0.33%	0.034%
24	14.093	1844	1854	1868	rBV	2609138	3606773	43.15%	4.429%
25	14.204	1868	1873	1877	rVB8	9448	16804	0.20%	0.021%
26	14.322	1890	1893	1896	rBV4	5126	8543	0.10%	0.010%
27	14.381	1896	1903	1913	rVV	2662248	4043448	48.38%	4.965%
28	14.463	1913	1917	1929	rVV5	20817	42357	0.51%	0.052%
29	14.563	1929	1934	1941	rVB2	10806	16882	0.20%	0.021%
30	14.687	1947	1955	1964	rBV	1351538	1984267	23.74%	2.436%
31	14.757	1964	1967	1973	rVB6	14202	23669	0.28%	0.029%
32	14.898	1983	1991	2012	rBV	847799	1793068	21.45%	2.202%
33	15.093	2020	2024	2033	rVB6	28436	52155	0.62%	0.064%
34	15.504	2090	2094	2102	rVB7	8024	14209	0.17%	0.017%
35	15.592	2103	2109	2112	rBV8	6077	10191	0.12%	0.013%
36	15.681	2117	2124	2132	rVV	3631883	5192136	62.12%	6.375%

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
 Data File : BP014225.D
 Acq On : 22 Mar 2023 19:54
 Operator : CG/JU
 Sample : 01977-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EYG88

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-BP030723.M
 Title : SVOA CALIBRATION

37	15.745	2133	2135	2139	rVV5	12932	19756	0.24%	0.024%
38	15.804	2139	2145	2158	rVV	1366841	1944060	23.26%	2.387%
39	15.922	2160	2165	2170	rVB4	19392	32492	0.39%	0.040%
40	15.998	2175	2178	2181	rVV5	12077	18888	0.23%	0.023%
41	16.045	2181	2186	2195	rVB	259872	351858	4.21%	0.432%
42	16.210	2210	2214	2218	rBV7	6036	9718	0.12%	0.012%
43	16.387	2240	2244	2252	rVB10	6445	13493	0.16%	0.017%
44	16.457	2252	2256	2264	rBV10	6939	15805	0.19%	0.019%
45	16.528	2265	2268	2273	rBV3	17620	27724	0.33%	0.034%
46	16.698	2292	2297	2304	rBV7	34033	64715	0.77%	0.079%
47	16.828	2314	2319	2338	rBV	269559	424197	5.08%	0.521%
48	17.045	2352	2356	2361	rBV8	4802	8585	0.10%	0.011%
49	17.116	2366	2368	2373	rVB5	7691	10346	0.12%	0.013%
50	17.187	2375	2380	2385	rBV8	25442	51950	0.62%	0.064%
51	17.239	2385	2389	2396	rVB4	34309	63623	0.76%	0.078%
52	17.328	2400	2404	2411	rBV5	32332	60453	0.72%	0.074%
53	17.392	2411	2415	2418	rBV	94249	128483	1.54%	0.158%
54	17.445	2418	2424	2435	rVV	1803523	2640679	31.59%	3.242%
55	17.545	2435	2441	2447	rVV	4119916	5774786	69.09%	7.091%
56	17.598	2447	2450	2466	rVB	219335	386339	4.62%	0.474%
57	17.804	2482	2485	2493	rVB8	22029	42389	0.51%	0.052%
58	17.934	2504	2507	2512	rVB6	18885	25227	0.30%	0.031%
59	17.981	2513	2515	2519	rBV5	9066	12641	0.15%	0.016%
60	18.057	2525	2528	2531	rBV4	19201	25130	0.30%	0.031%
61	18.186	2545	2550	2560	rBV	500907	778922	9.32%	0.956%
62	18.310	2566	2571	2581	rBV	2323561	3088661	36.95%	3.792%
63	18.598	2615	2620	2623	rBV7	24995	49111	0.59%	0.060%
64	18.657	2624	2630	2635	rVV6	69355	161556	1.93%	0.198%
65	18.710	2635	2639	2647	rVV	104422	130242	1.56%	0.160%
66	18.781	2649	2651	2655	rVV5	18041	22430	0.27%	0.028%
67	18.834	2656	2660	2672	rVB5	67371	136129	1.63%	0.167%
68	18.951	2676	2680	2687	rBV5	42160	74242	0.89%	0.091%
69	19.063	2697	2699	2703	rBV5	17748	20565	0.25%	0.025%
70	19.128	2707	2710	2717	rVV7	37626	52008	0.62%	0.064%
71	19.351	2744	2748	2752	rBV2	70967	103390	1.24%	0.127%
72	19.416	2752	2759	2771	rVV	362854	744101	8.90%	0.914%
73	19.533	2775	2779	2789	rBV2	314380	468183	5.60%	0.575%
74	19.769	2813	2819	2825	rBV	5070686	6939593	83.03%	8.521%
75	20.263	2900	2903	2907	rVB	70397	73696	0.88%	0.090%
76	20.745	2982	2985	2989	rBV2	100957	109588	1.31%	0.135%

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
 Data File : BP014225.D
 Acq On : 22 Mar 2023 19:54
 Operator : CG/JU
 Sample : 01977-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EYG88

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-BP030723.M

Title : SVOA CALIBRATION

77	21.198	3058	3062	3068	rBV	1129606	1371593	16.41%	1.684%
78	21.404	3094	3097	3103	rBV	212557	234827	2.81%	0.288%
79	21.527	3113	3118	3128	rBV	2197081	3007835	35.99%	3.693%
80	21.651	3136	3139	3142	rVB	106122	117646	1.41%	0.144%
81	22.798	3328	3334	3343	rBV	5592642	8358312	100.00%	10.263%
82	23.886	3512	3519	3530	rVB2	3283477	6920320	82.80%	8.497%
83	24.051	3541	3547	3557	rVB2	1364490	2896087	34.65%	3.556%

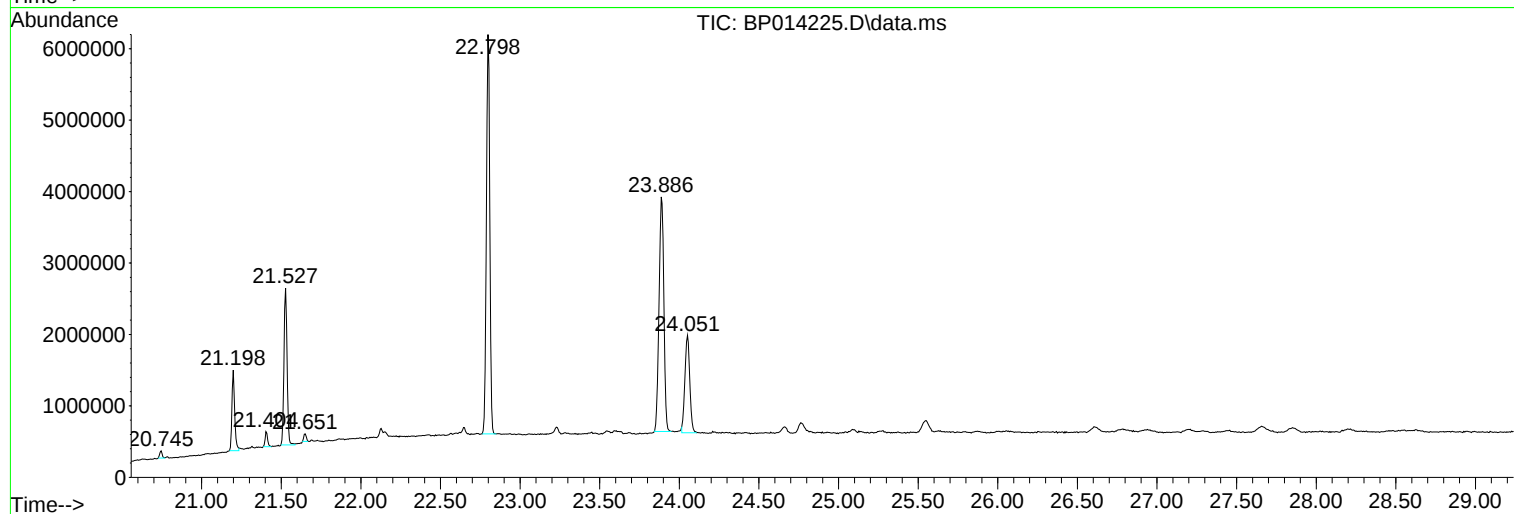
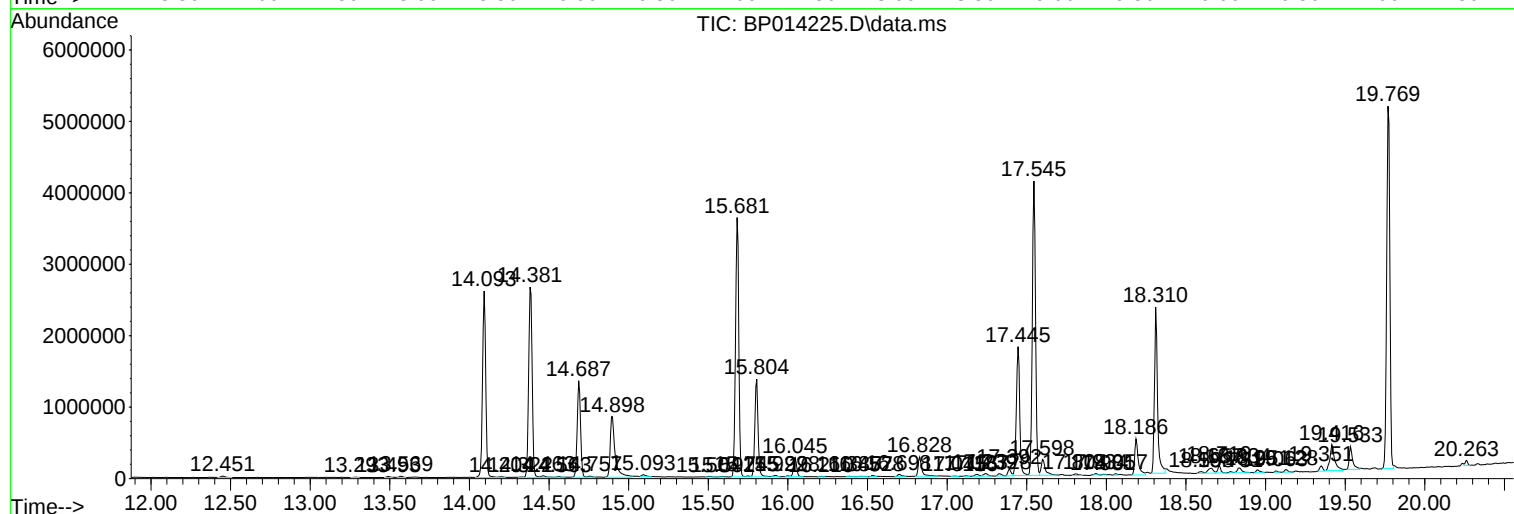
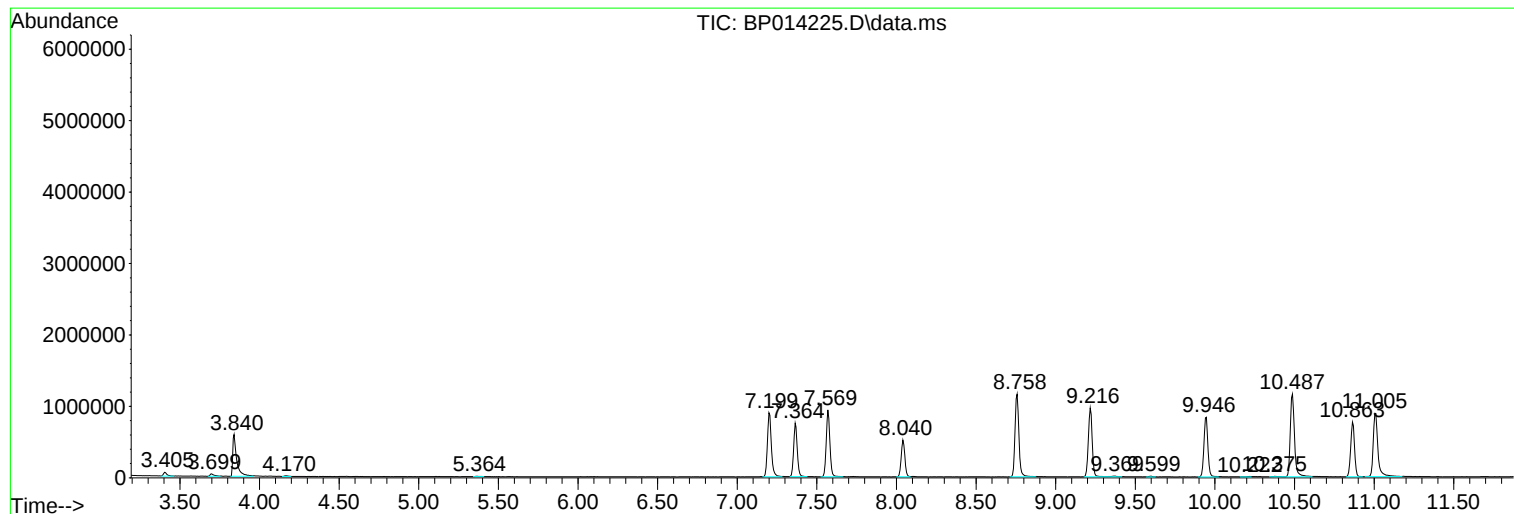
Sum of corrected areas: 81443213

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
Data File : BP014225.D
Acq On : 22 Mar 2023 19:54
Operator : CG/JU
Sample : 01977-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EYG88

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
Data File : BP014225.D
Acq On : 22 Mar 2023 19:54
Operator : CG/JU
Sample : 01977-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EYG88

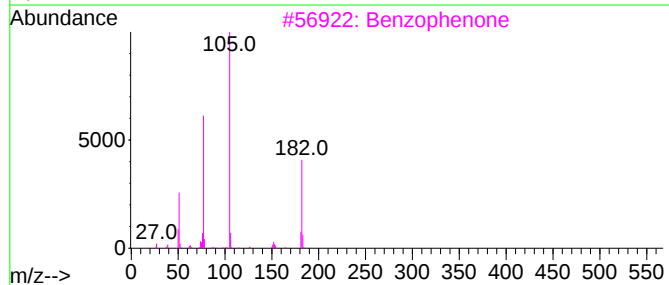
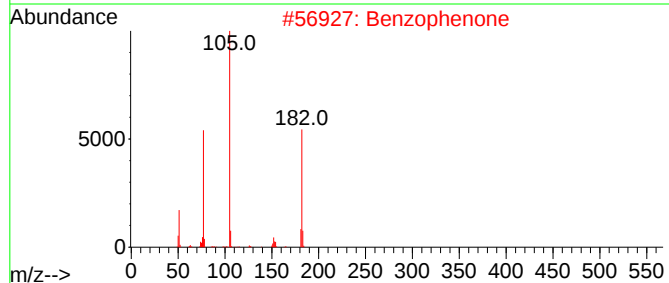
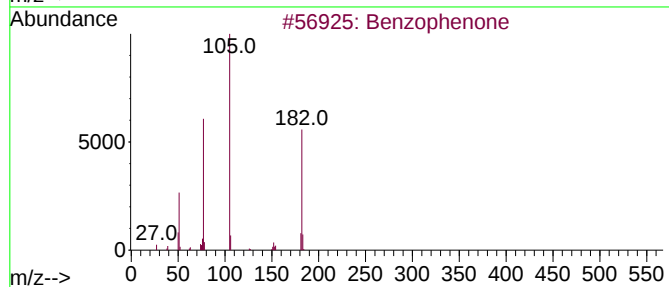
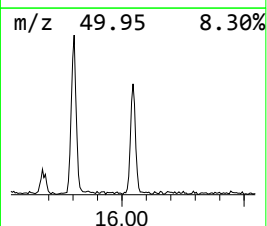
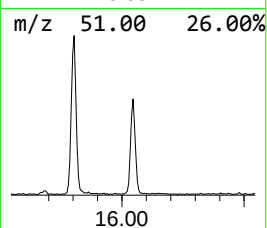
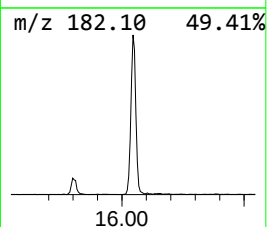
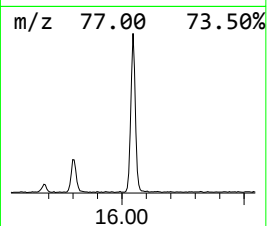
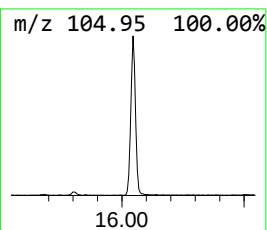
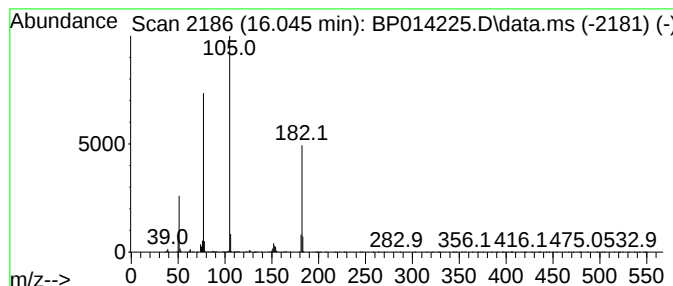
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.045	3.55 ng/ul	351858	Acenaphthene-d10	14.687

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	94
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
Data File : BP014225.D
Acq On : 22 Mar 2023 19:54
Operator : CG/JU
Sample : 01977-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EYG88

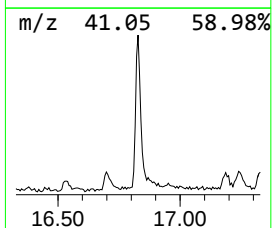
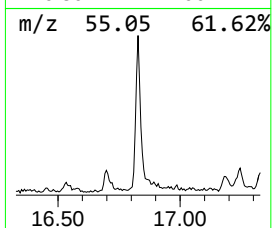
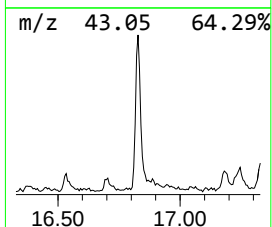
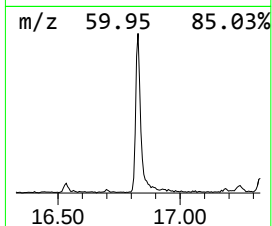
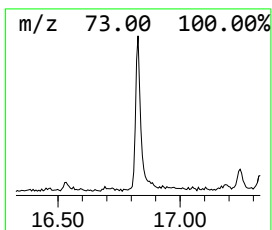
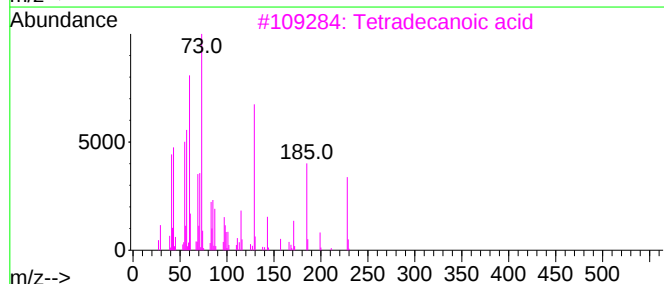
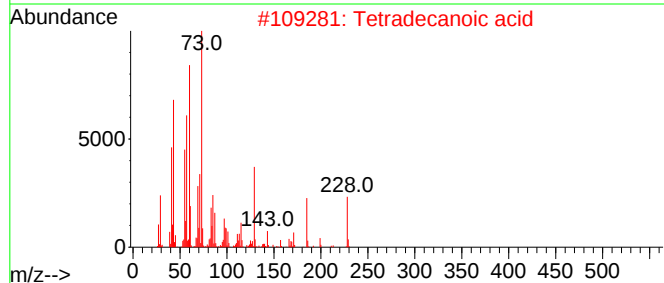
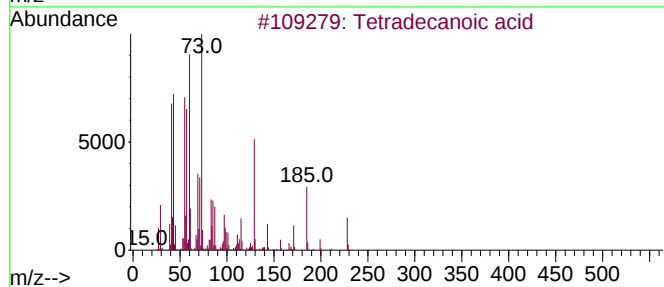
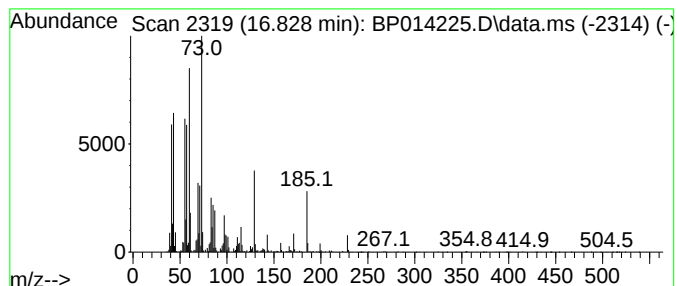
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Tetradecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.828	3.21 ng/ul	424197	Phenanthrene-d10	17.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecanoic acid	228	C14H28O2	000544-63-8	98
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	90
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	90
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	87
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
Data File : BP014225.D
Acq On : 22 Mar 2023 19:54
Operator : CG/JU
Sample : 01977-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EYG88

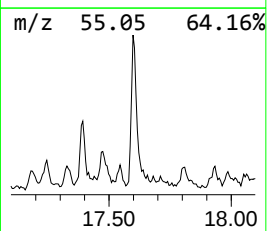
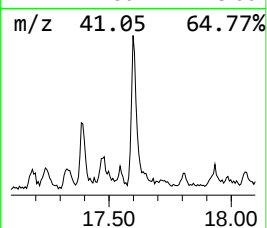
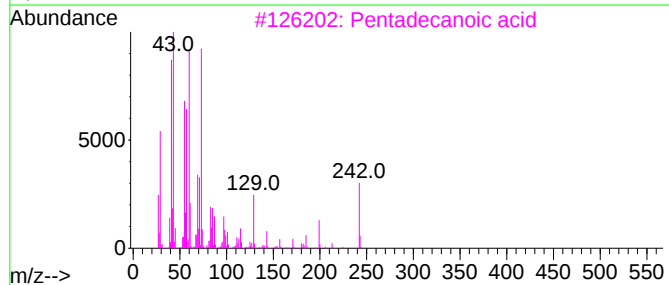
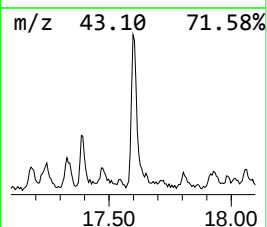
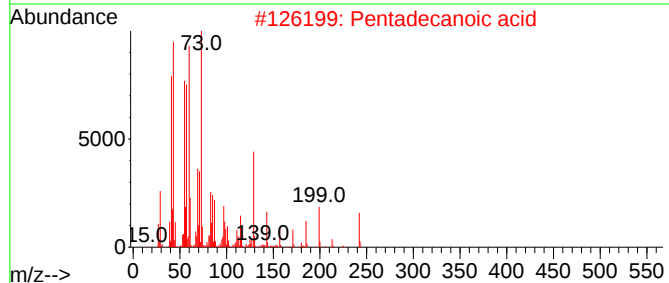
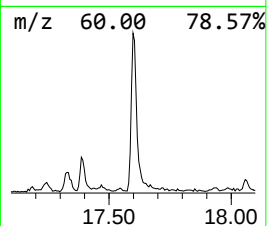
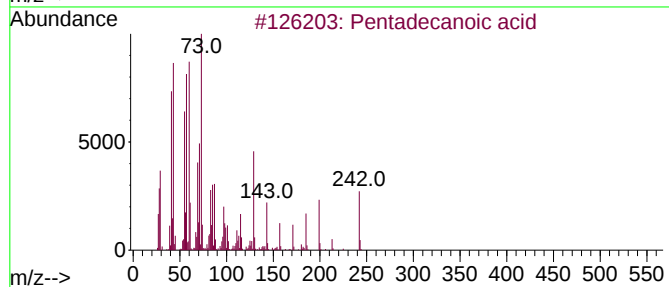
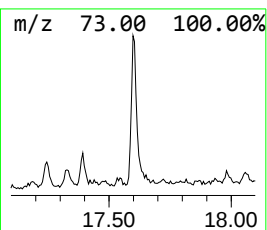
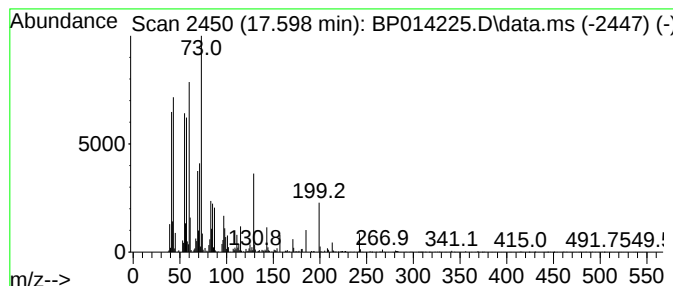
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Pentadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.598	2.93 ng/ul	386339	Phenanthrene-d10	17.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecanoic acid	242	C15H30O2	001002-84-2	98
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	98
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	96
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	86
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
Data File : BP014225.D
Acq On : 22 Mar 2023 19:54
Operator : CG/JU
Sample : 01977-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EYG88

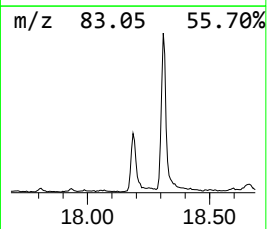
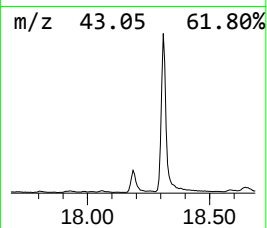
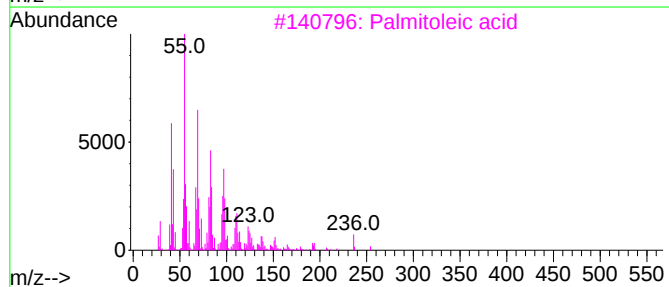
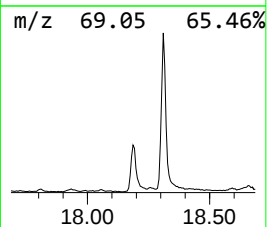
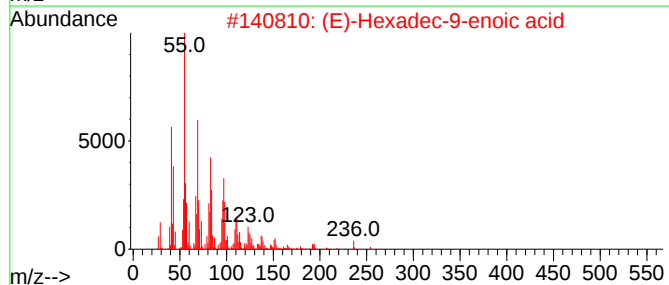
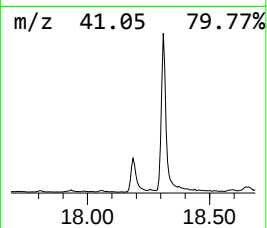
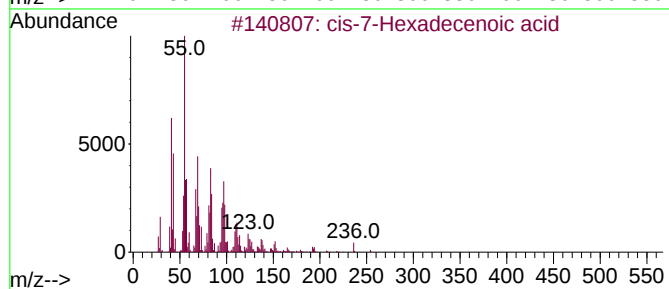
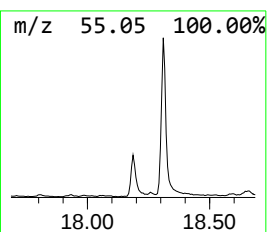
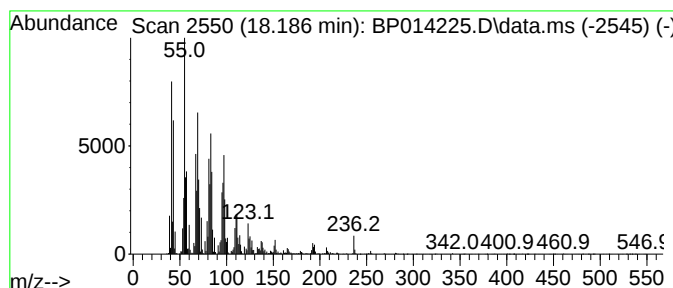
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 cis-7-Hexadecenoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.186	5.90 ng/ul	778922	Phenanthrene-d10	17.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cis-7-Hexadecenoic acid	254	C16H30O2	002416-19-5	99
2			(E)-Hexadec-9-enoic acid	254	C16H30O2	010030-73-6	99
3			Palmitoleic acid	254	C16H30O2	000373-49-9	97
4			cis-7-Hexadecenoic acid	254	C16H30O2	002416-19-5	96
5			Palmitoleic acid	254	C16H30O2	000373-49-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
Data File : BP014225.D
Acq On : 22 Mar 2023 19:54
Operator : CG/JU
Sample : 01977-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

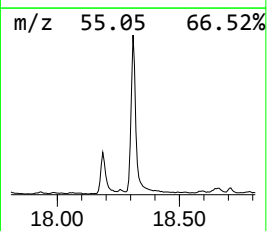
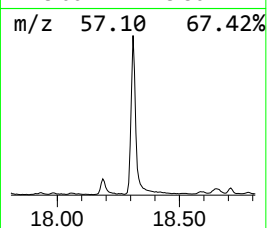
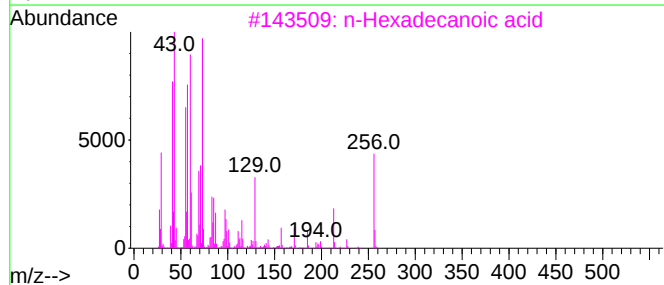
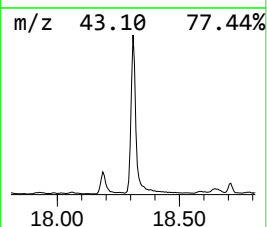
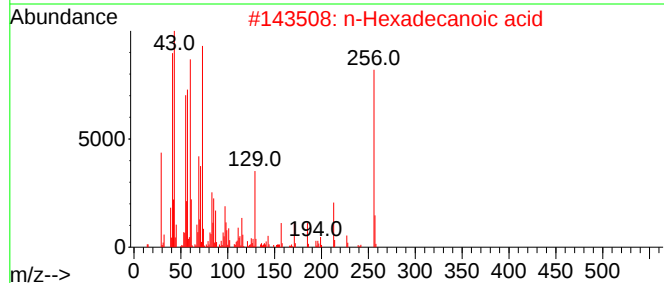
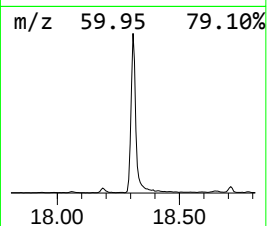
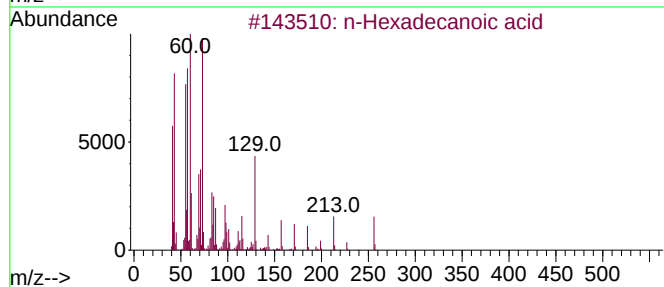
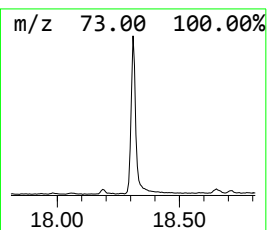
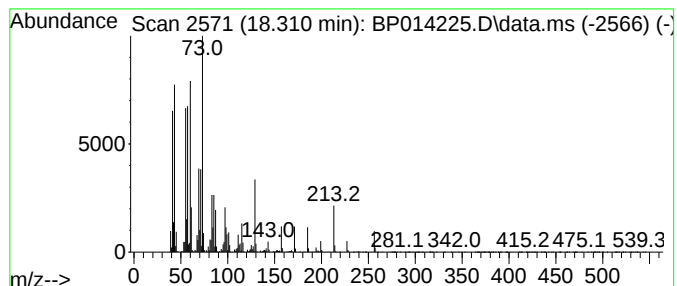
Instrument :
BNA_P
ClientSampleId :
EYG88

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
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.	
18.310	23.39 ng/ul	3088660	Phenanthrene-d10	17.445	
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	98
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
Data File : BP014225.D
Acq On : 22 Mar 2023 19:54
Operator : CG/JU
Sample : 01977-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EYG88

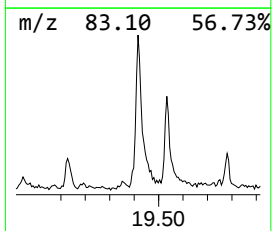
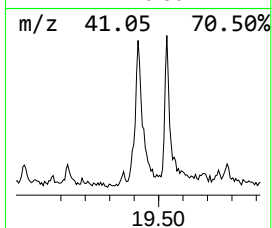
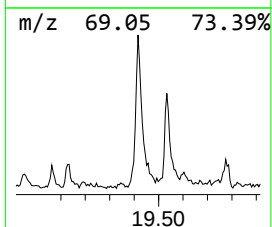
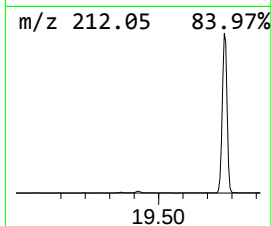
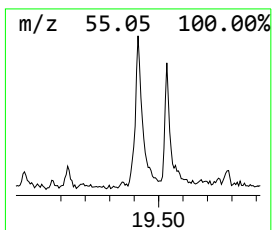
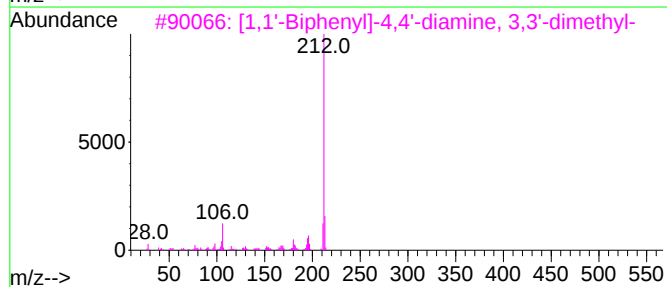
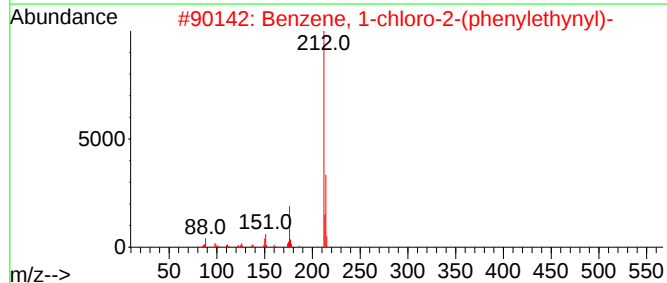
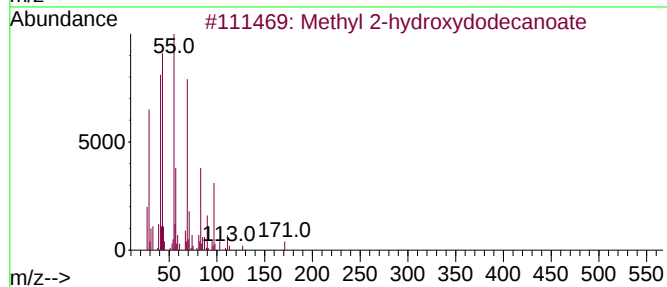
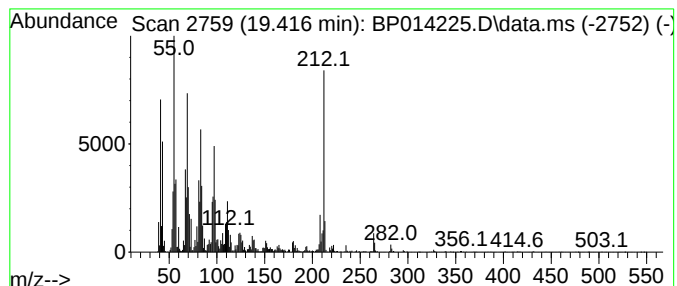
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.416	5.64 ng/ul	744101	Phenanthrene-d10	17.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 2-hydroxydodecanoate	230	C13H26O3	051067-85-7	35
2			Benzene, 1-chloro-2-(phenylethyn...	212	C14H9Cl	010271-57-5	30
3			[1,1'-Biphenyl]-4,4'-diamine, 3,...	212	C14H16N2	000119-93-7	25
4			Benzoic acid, 3,4,5-trimethoxy-	212	C10H12O5	000118-41-2	15
5			4-Amino-1,8-naphthalimide	212	C12H8N2O2	001742-95-6	15



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
Data File : BP014225.D
Acq On : 22 Mar 2023 19:54
Operator : CG/JU
Sample : 01977-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EYG88

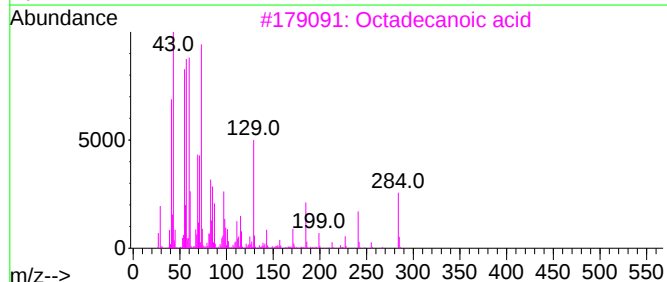
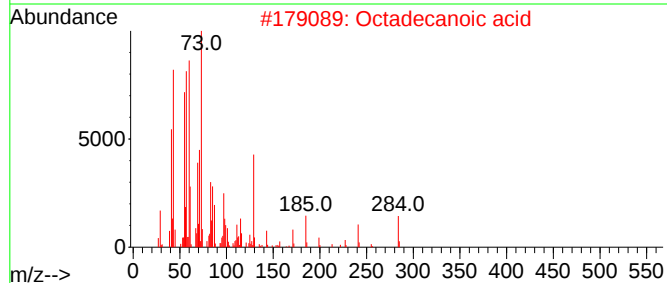
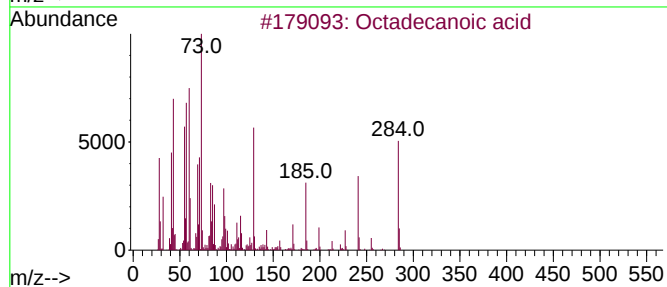
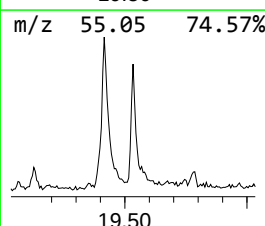
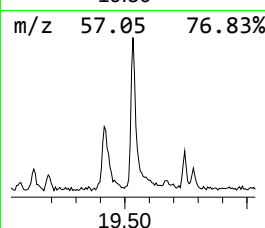
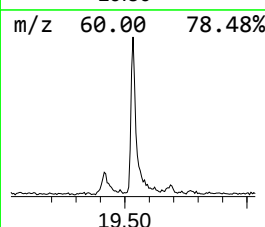
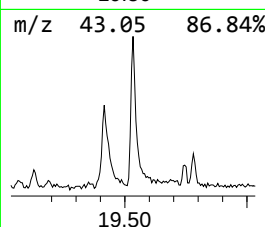
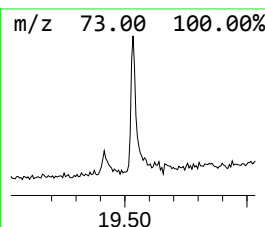
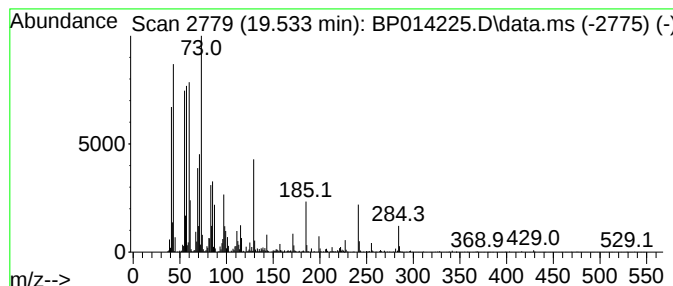
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 Octadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.534	3.11 ng/ul	468183	Chrysene-d12	21.528

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	99
4			Octadecanoic acid	284	C18H36O2	000057-11-4	98
5			Octadecanoic acid	284	C18H36O2	000057-11-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
Data File : BP014225.D
Acq On : 22 Mar 2023 19:54
Operator : CG/JU
Sample : 01977-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EYG88

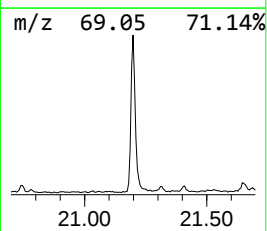
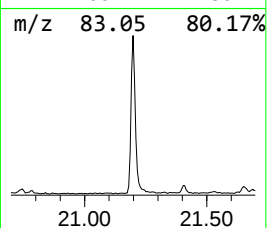
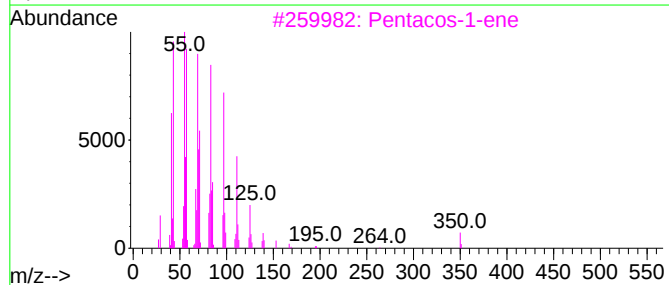
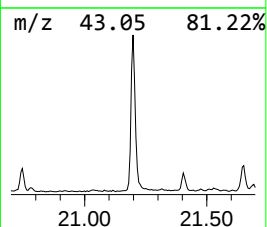
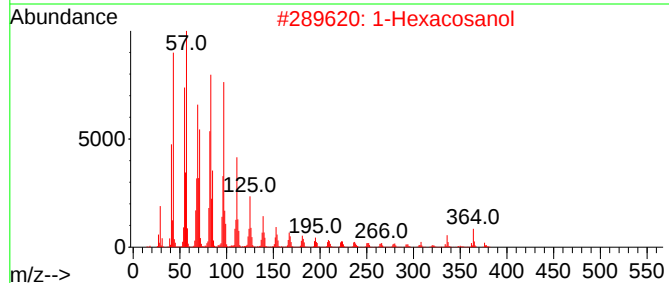
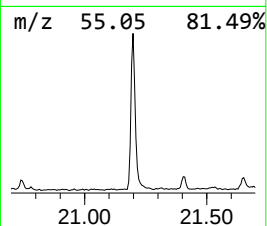
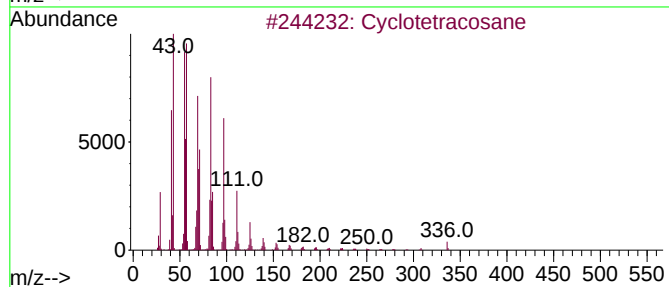
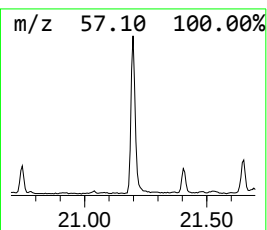
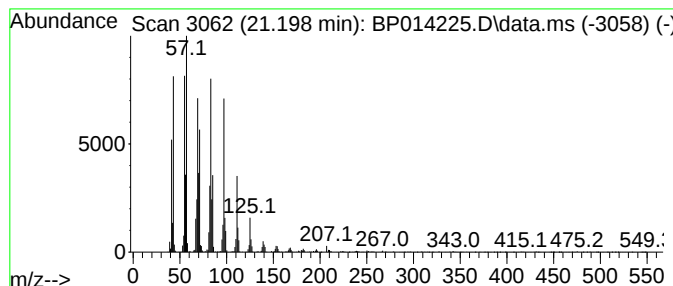
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Cyclic21.198 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.198	9.12 ng/ul	1371590	Chrysene-d12	21.528

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclotetracosane	336	C24H48	000297-03-0	94
2			1-Hexacosanol	382	C26H54O	000506-52-5	94
3			Pentacos-1-ene	350	C25H50	016980-85-1	94
4			Ethanol, 2-(tetradecyloxy)-	258	C16H34O2	002136-70-1	93
5			Pentafluoropropionic acid, hexad...	388	C19H33F5O2	006222-07-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
Data File : BP014225.D
Acq On : 22 Mar 2023 19:54
Operator : CG/JU
Sample : 01977-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EYG88

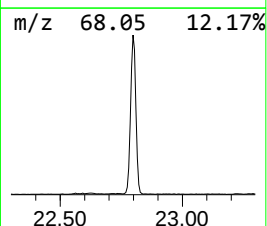
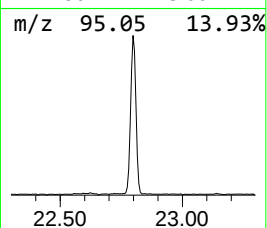
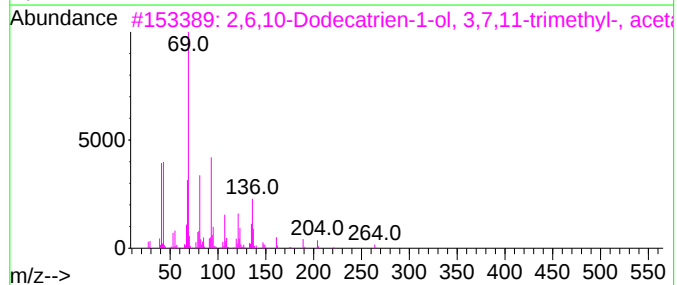
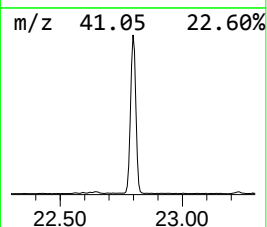
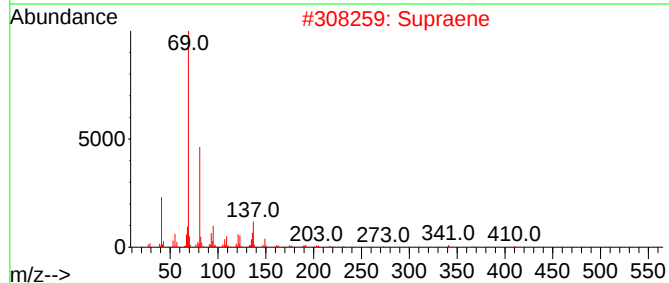
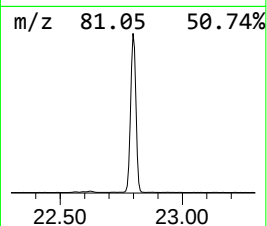
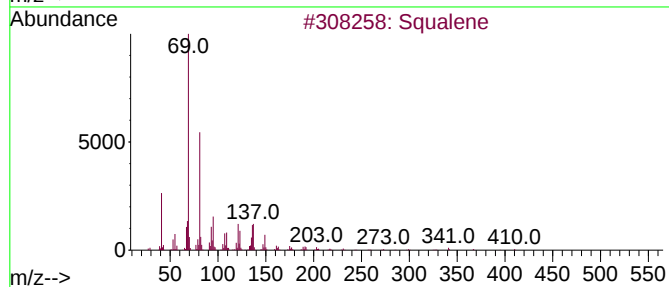
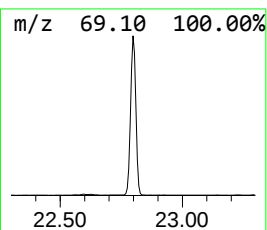
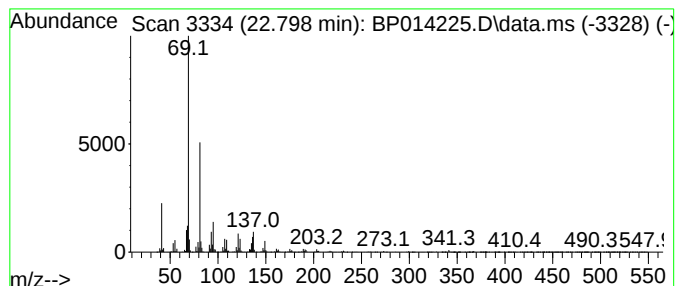
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 Squalene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.798	57.72 ng/ul	8358310	Perylene-d12	24.051

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Squalene	410	C30H50	000111-02-4	94
2			Supraene	410	C30H50	007683-64-9	91
3			2,6,10-Dodecatrien-1-ol, 3,7,11-...	264	C17H28O2	004128-17-0	64
4			3,7-Dimethyl-1-phenylsulfonyl-2,...	278	C16H22O2S	1000432-27-5	59
5			2,3,5-Trimethyl-2,3,5-hexanetric...	203	C12H17N3	003974-79-6	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
Data File : BP014225.D
Acq On : 22 Mar 2023 19:54
Operator : CG/JU
Sample : 01977-02
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EYG88

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.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
Benzophenone	16.045	3.5	ng/ul	351858	3	14.687	1984270	20.0
Tetradecanoic acid	16.828	3.2	ng/ul	424197	4	17.445	2640680	20.0
Pentadecanoic acid	17.598	2.9	ng/ul	386339	4	17.445	2640680	20.0
cis-7-Hexadecen...	18.186	5.9	ng/ul	778922	4	17.445	2640680	20.0
n-Hexadecanoic ...	18.310	23.4	ng/ul	3088660	4	17.445	2640680	20.0
unknown-01	19.416	5.6	ng/ul	744101	4	17.445	2640680	20.0
Octadecanoic acid	19.534	3.1	ng/ul	468183	5	21.528	3007840	20.0
(DEL) Alkane: C...	21.198	9.1	ng/ul	1371590	5	21.528	3007840	20.0
Squalene	22.798	57.7	ng/ul	8358310	6	24.051	2896090	20.0