

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
 Data File : BP015865.D
 Acq On : 01 Jul 2023 03:02
 Operator : MA/JU
 Sample : 03161-15
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFX1

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-BP062623.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP015865.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.375	29	32	38	rVB	30428	43077	0.99%	0.079%
2	3.822	105	108	110	rBV	17915	20567	0.47%	0.038%
3	3.846	110	112	122	rVV	71141	139321	3.21%	0.255%
4	3.928	122	126	132	rVV	85116	131540	3.03%	0.241%
5	4.005	136	139	141	rVV	15398	20396	0.47%	0.037%
6	4.046	142	146	154	rVB	38393	60668	1.40%	0.111%
7	4.146	160	163	173	rVB	19402	38396	0.88%	0.070%
8	4.728	259	262	267	rVB6	11613	15168	0.35%	0.028%
9	5.093	321	324	329	rBV	10786	16022	0.37%	0.029%
10	6.510	561	565	569	rBV7	3010	6509	0.15%	0.012%
11	7.246	683	690	707	rBV2	215472	674718	15.53%	1.235%
12	7.381	707	713	731	rVB	350047	673685	15.51%	1.233%
13	7.587	742	748	769	rBV	372460	800147	18.42%	1.465%
14	8.051	820	827	847	rBV	910470	1788473	41.17%	3.274%
15	8.822	951	958	972	rBV3	61336	252600	5.81%	0.462%
16	9.251	1024	1031	1052	rBV	342282	828978	19.08%	1.518%
17	9.592	1083	1089	1092	rVB4	7566	13323	0.31%	0.024%
18	9.981	1147	1155	1178	rBV2	202787	711250	16.37%	1.302%
19	10.563	1247	1254	1285	rBV	205981	862312	19.85%	1.579%
20	10.881	1299	1308	1328	rBV	1274048	2768059	63.71%	5.067%
21	11.092	1336	1344	1368	rBV	74267	329499	7.58%	0.603%
22	11.269	1372	1374	1379	rVB6	3959	6150	0.14%	0.011%
23	11.892	1478	1480	1486	rVB7	3661	5615	0.13%	0.010%
24	12.081	1509	1512	1517	rVV6	6111	9045	0.21%	0.017%
25	12.128	1517	1520	1524	rVB6	3312	4828	0.11%	0.009%
26	12.281	1544	1546	1551	rVV6	3890	5796	0.13%	0.011%
27	12.410	1564	1568	1571	rBV6	4327	6033	0.14%	0.011%
28	12.516	1580	1586	1593	rVB10	6434	16522	0.38%	0.030%
29	13.216	1699	1705	1708	rBV8	3959	7476	0.17%	0.014%
30	13.504	1750	1754	1758	rBV6	6826	9706	0.22%	0.018%
31	13.580	1761	1767	1775	rBV5	16446	34580	0.80%	0.063%
32	13.675	1780	1783	1788	rVB7	3538	6399	0.15%	0.012%
33	13.957	1828	1831	1836	rBV7	3285	5772	0.13%	0.011%
34	14.069	1845	1850	1851	rBV5	3536	5080	0.12%	0.009%
35	14.116	1851	1858	1883	rBV	885731	1670853	38.46%	3.059%
36	14.398	1899	1906	1929	rBV	1202479	1965896	45.25%	3.599%

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
 Data File : BP015865.D
 Acq On : 01 Jul 2023 03:02
 Operator : MA/JU
 Sample : 03161-15
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFX1

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-BP062623.MA.M
 Title : SVOA CALIBRATION

37	14.569	1932	1935	1941	rVB6	9985	14383	0.33%	0.026%
38	14.704	1950	1958	1976	rBV	2225552	3674163	84.57%	6.726%
39	15.027	2006	2013	2026	rBV4	61271	283149	6.52%	0.518%
40	15.469	2084	2088	2092	rVV6	6298	9686	0.22%	0.018%
41	15.522	2094	2097	2103	rVB7	15608	22987	0.53%	0.042%
42	15.704	2121	2128	2143	rBV	1309015	2415882	55.61%	4.423%
43	15.880	2153	2158	2179	rBV3	121525	405538	9.33%	0.742%
44	16.074	2185	2191	2204	rBV	316787	559010	12.87%	1.023%
45	16.386	2240	2244	2255	rBV10	14927	38552	0.89%	0.071%
46	16.557	2269	2273	2276	rBV6	11972	16949	0.39%	0.031%
47	16.633	2284	2286	2293	rVB6	15372	20272	0.47%	0.037%
48	16.851	2319	2323	2325	rBV5	13228	21367	0.49%	0.039%
49	16.892	2327	2330	2337	rVB8	22043	42497	0.98%	0.078%
50	16.998	2345	2348	2356	rVB5	20839	29436	0.68%	0.054%
51	17.180	2376	2379	2381	rBV3	14859	18500	0.43%	0.034%
52	17.339	2402	2406	2413	rBV5	34623	65166	1.50%	0.119%
53	17.404	2414	2417	2421	rVB5	19437	24210	0.56%	0.044%
54	17.463	2421	2427	2440	rBV	2541021	4166918	95.91%	7.628%
55	17.569	2440	2445	2465	rVB2	1101488	2136079	49.17%	3.910%
56	17.916	2502	2504	2509	rBV3	16251	20497	0.47%	0.038%
57	18.098	2532	2535	2539	rVB3	28817	34383	0.79%	0.063%
58	18.204	2551	2553	2558	rVB3	32591	34304	0.79%	0.063%
59	18.263	2558	2563	2568	rBV	449123	648812	14.93%	1.188%
60	18.316	2568	2572	2582	rVV	1568207	2098258	48.30%	3.841%
61	19.192	2718	2721	2724	rBV	44668	47522	1.09%	0.087%
62	19.404	2754	2757	2759	rBV2	124757	159537	3.67%	0.292%
63	19.427	2759	2761	2762	rVV	144542	131836	3.03%	0.241%
64	19.451	2763	2765	2776	rVB4	208597	373998	8.61%	0.685%
65	19.539	2776	2780	2786	rBV	361572	572514	13.18%	1.048%
66	19.792	2819	2823	2851	rVB	2010935	4239349	97.58%	7.761%
67	20.268	2901	2904	2910	rVB2	143913	173461	3.99%	0.318%
68	20.757	2982	2987	2991	rBV5	346725	853490	19.64%	1.562%
69	21.221	3060	3066	3073	rBV2	682246	1230660	28.33%	2.253%
70	21.415	3096	3099	3103	rBV	151515	168862	3.89%	0.309%
71	21.557	3118	3123	3133	rBV	1905077	3130022	72.04%	5.730%
72	22.127	3217	3220	3224	rVB	387219	446350	10.27%	0.817%
73	23.233	3403	3408	3415	rVB	723514	1191017	27.41%	2.180%
74	24.098	3549	3555	3573	rVB	1427761	3748997	86.29%	6.863%
75	24.662	3645	3651	3664	rVB	1396014	3056669	70.36%	5.596%
76	27.262	4087	4093	4106	rVB4	1386449	4344573	100.00%	7.954%

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015865.D
Acq On : 01 Jul 2023 03:02
Operator : MA/JU
Sample : 03161-15
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFX1

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-BP062623.MA.M
Title : SVOA CALIBRATION

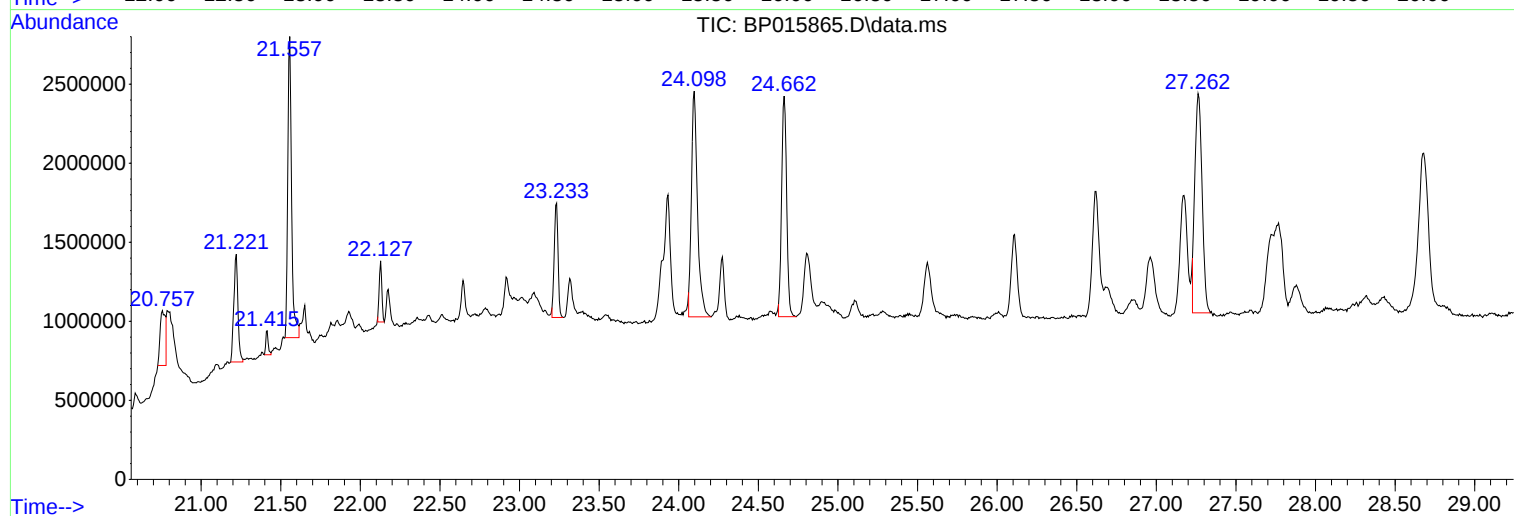
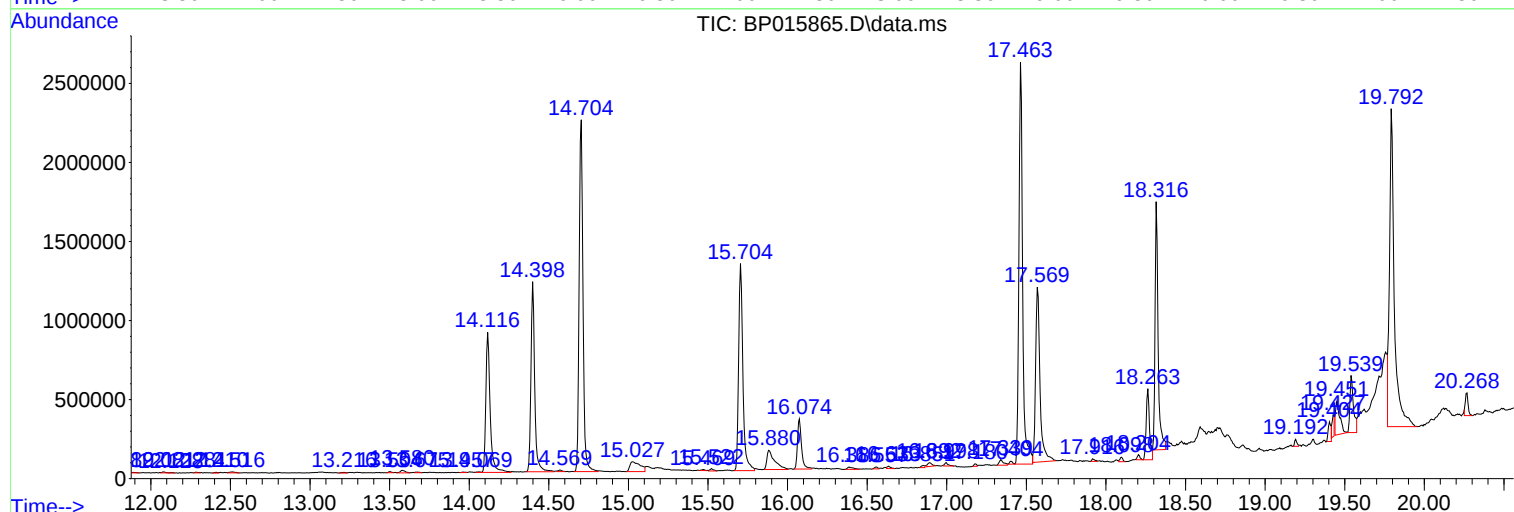
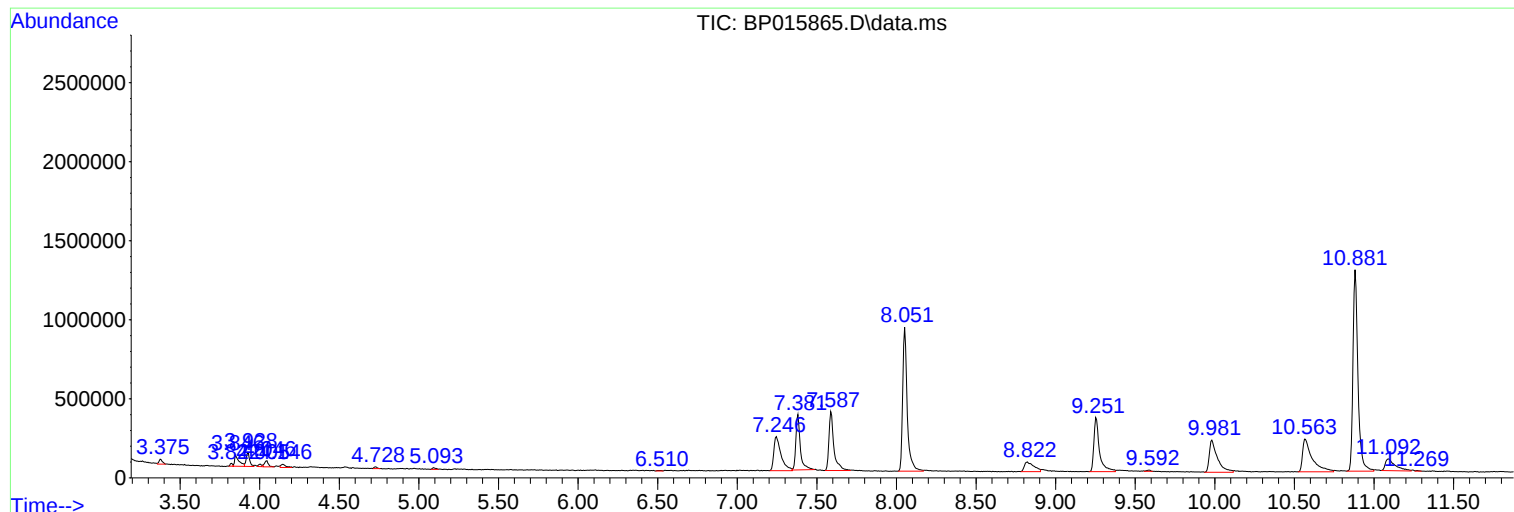
Sum of corrected areas: 54624314

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015865.D
Acq On : 01 Jul 2023 03:02
Operator : MA/JU
Sample : 03161-15
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFX1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.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\BP063023\
Data File : BP015865.D
Acq On : 01 Jul 2023 03:02
Operator : MA/JU
Sample : 03161-15
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFX1

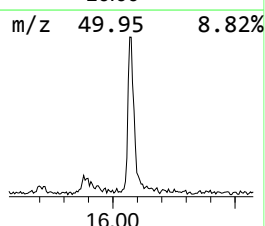
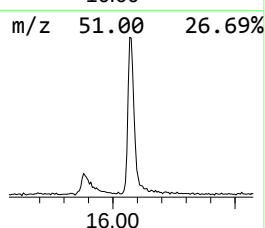
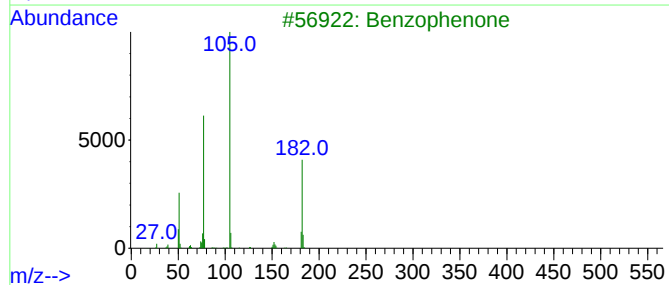
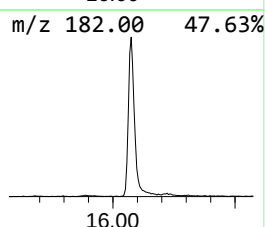
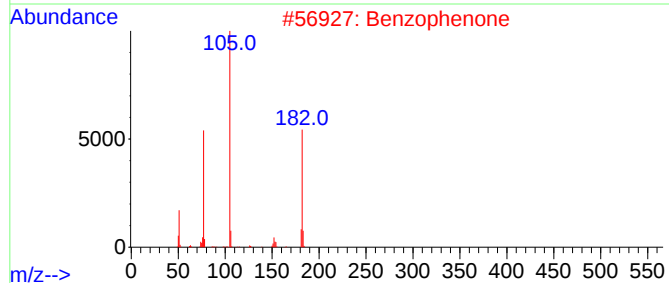
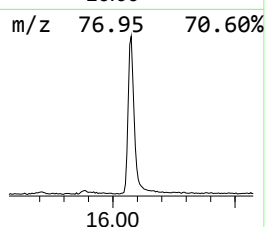
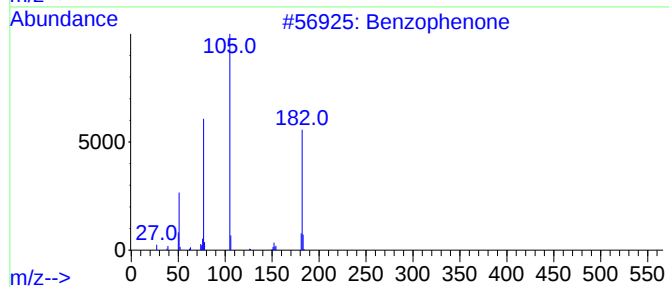
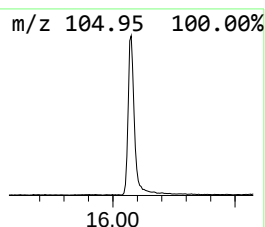
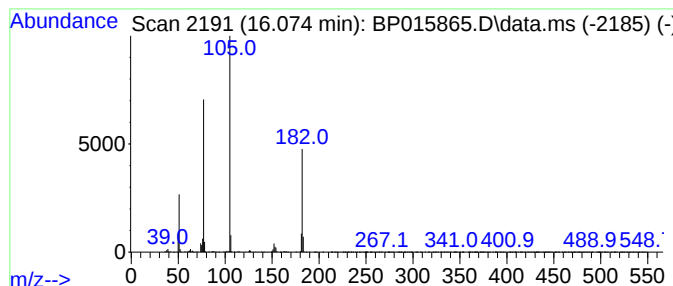
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.075	3.04 ng/ul	559010	Acenaphthene-d10	14.704

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
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	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015865.D
Acq On : 01 Jul 2023 03:02
Operator : MA/JU
Sample : 03161-15
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFX1

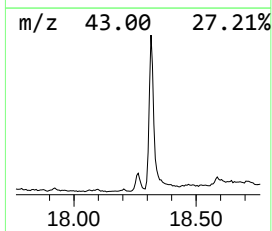
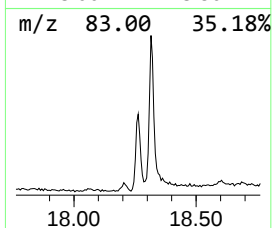
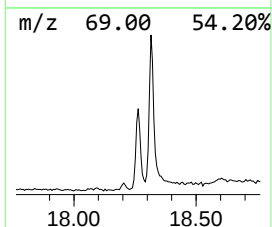
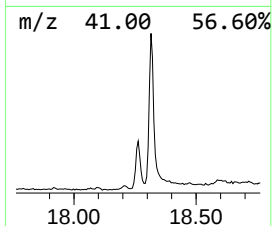
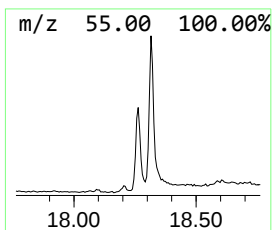
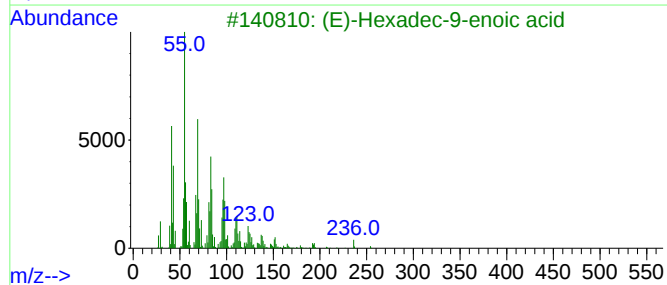
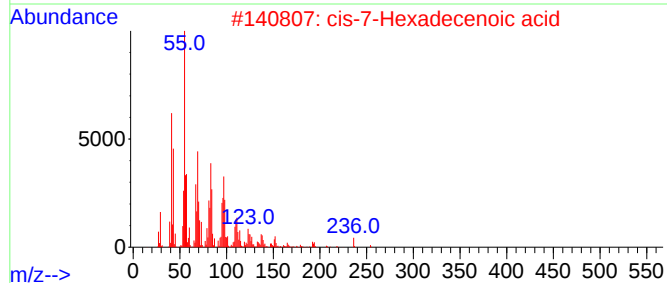
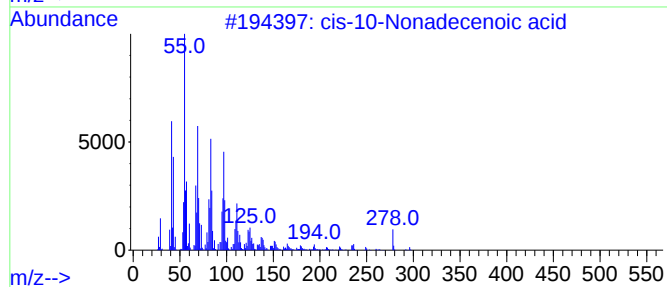
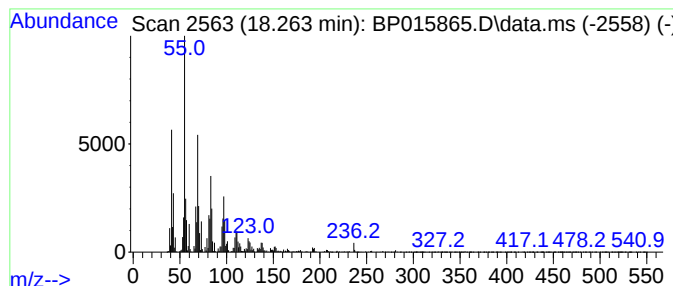
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 cis-10-Nonadecenoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.263	3.11 ng/ul	648812	Phenanthrene-d10	17.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cis-10-Nonadecenoic acid	296	C19H36O2	073033-09-7	99
2			cis-7-Hexadecenoic acid	254	C16H30O2	002416-19-5	97
3			(E)-Hexadec-9-enoic acid	254	C16H30O2	010030-73-6	96
4			Myristoleic acid	226	C14H26O2	000544-64-9	95
5			Myristoleic acid	226	C14H26O2	000544-64-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015865.D
Acq On : 01 Jul 2023 03:02
Operator : MA/JU
Sample : 03161-15
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFX1

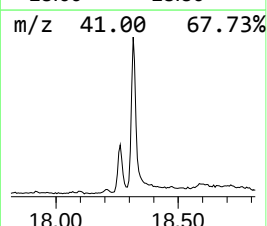
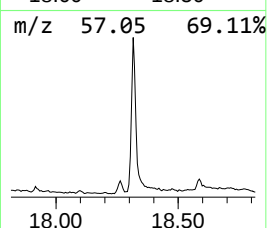
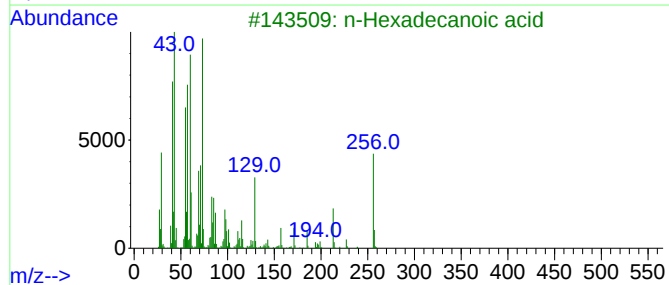
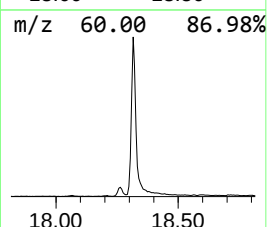
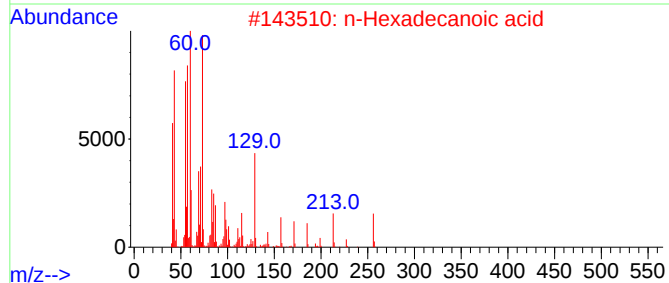
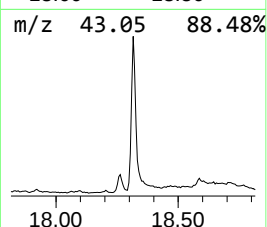
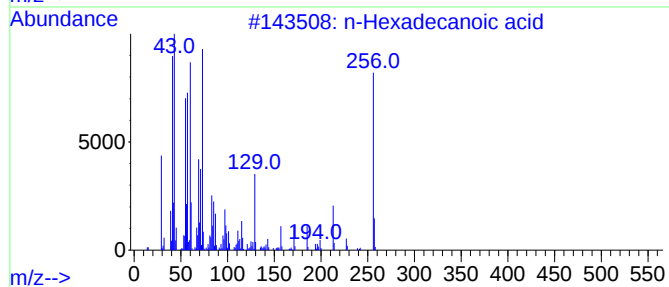
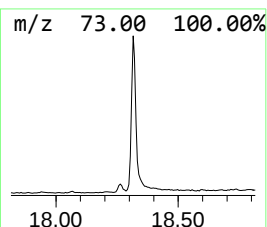
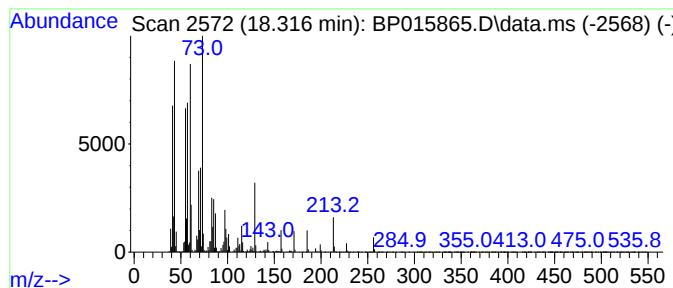
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.316	10.07 ng/ul	2098260	Phenanthrene-d10	17.463

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	97
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015865.D
Acq On : 01 Jul 2023 03:02
Operator : MA/JU
Sample : 03161-15
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFX1

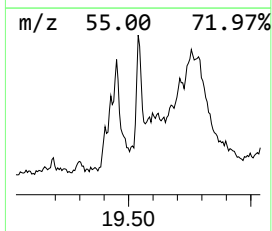
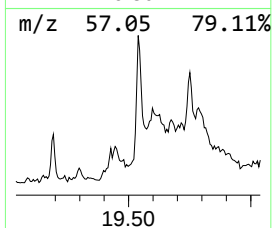
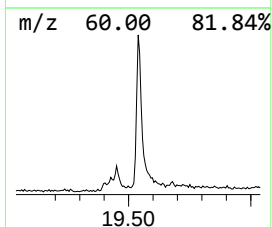
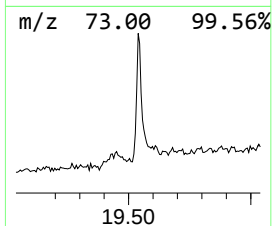
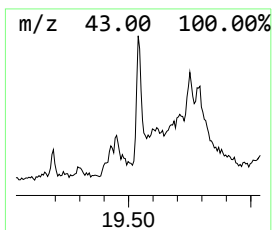
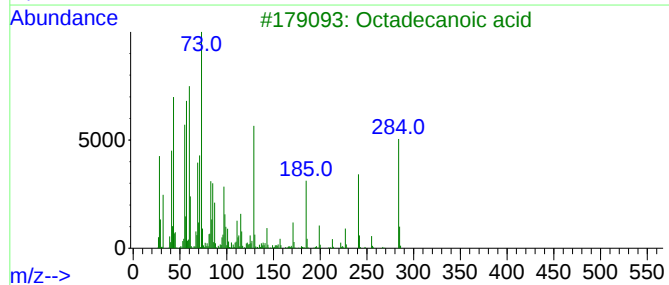
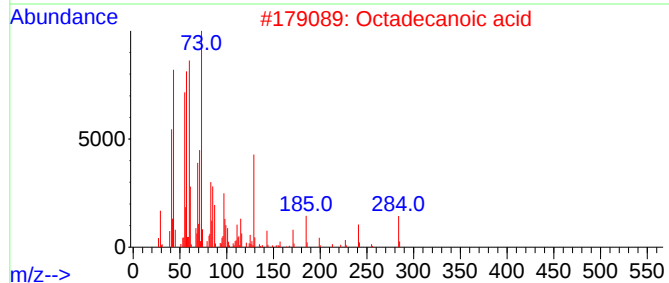
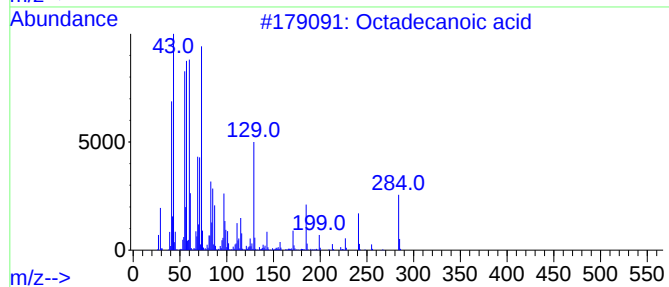
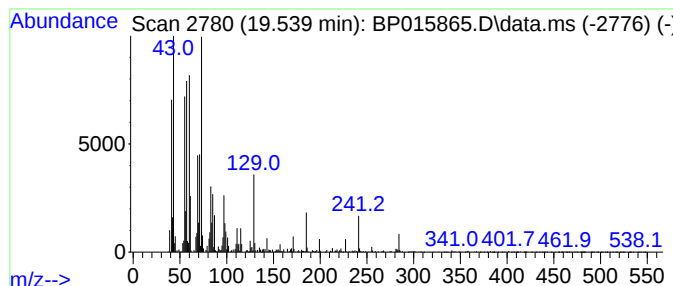
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.539	3.66 ng/ul	572514	Chrysene-d12	21.557

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	95
5			Octadecanoic acid	284	C18H36O2	000057-11-4	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015865.D
Acq On : 01 Jul 2023 03:02
Operator : MA/JU
Sample : 03161-15
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFX1

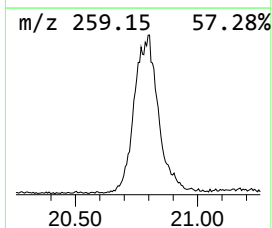
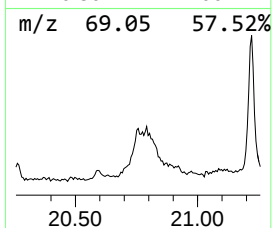
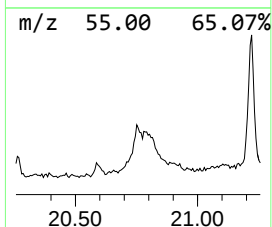
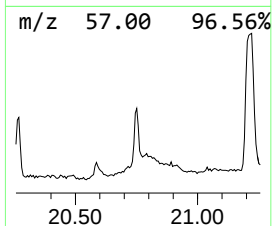
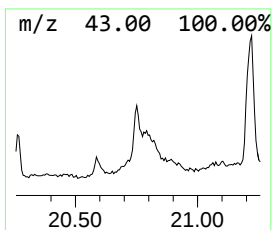
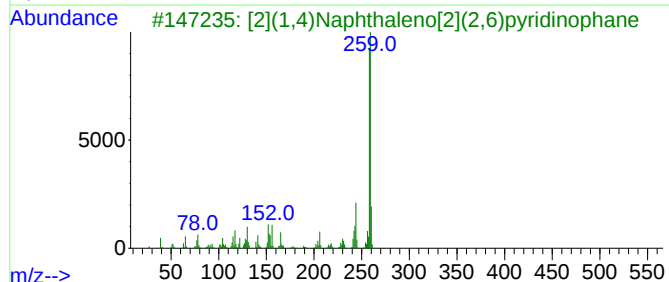
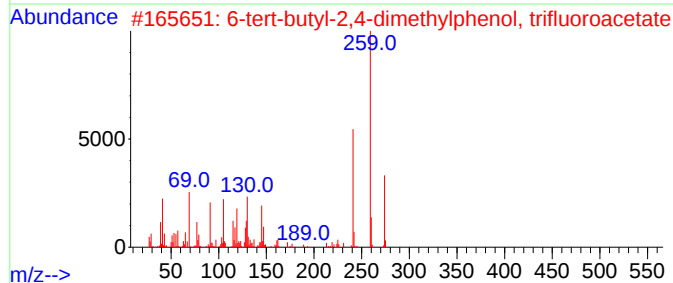
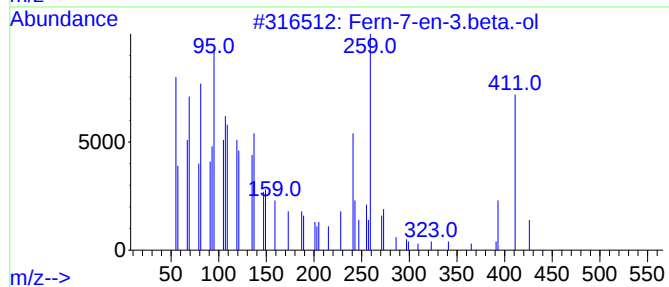
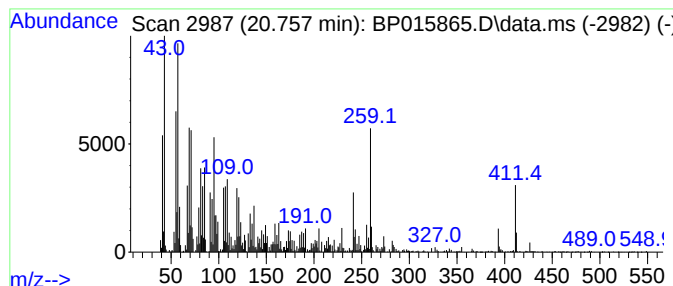
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Fern-7-en-3.beta.-ol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.757	5.45 ng/ul	853490	Chrysene-d12	21.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fern-7-en-3.beta.-ol	426	C30H50O	004966-00-1	89
2			6-tert-butyl-2,4-dimethylphenol,...	274	C14H17F3O2	1000376-41-4	18
3			[2](1,4)Naphthaleno[2](2,6)pyrid...	259	C19H17N	099677-63-1	15
4			Carbamic acid, (5-chloro-2,4-dim...	259	C11H14ClNO4	1000319-49-5	15
5			1-(4-Fluorophenyl)-3-(3-nitrophe...	260	C12H9FN4O2	1000410-81-9	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015865.D
Acq On : 01 Jul 2023 03:02
Operator : MA/JU
Sample : 03161-15
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFX1

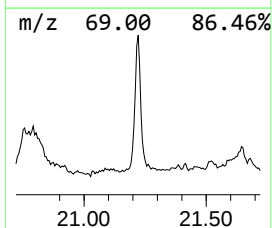
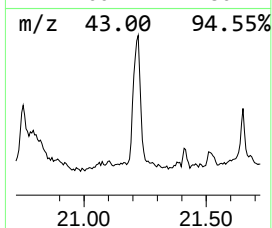
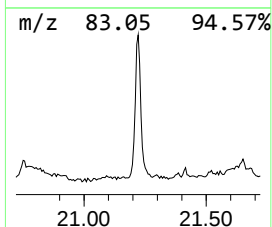
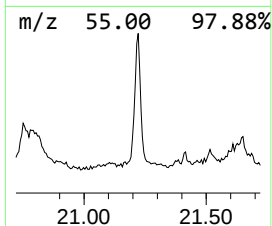
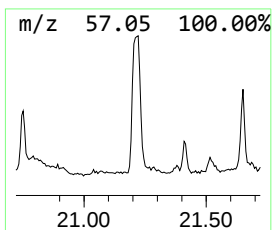
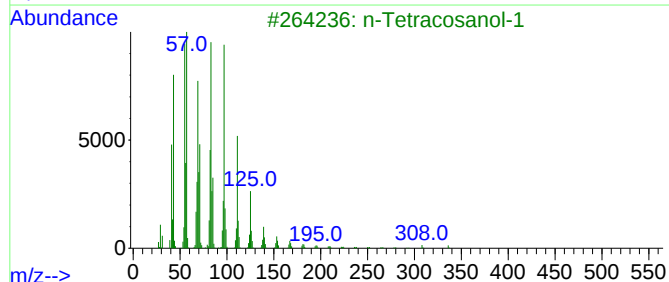
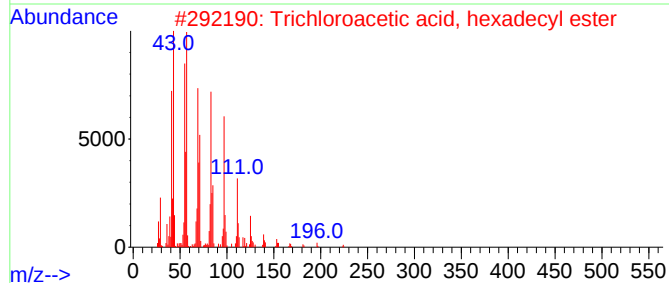
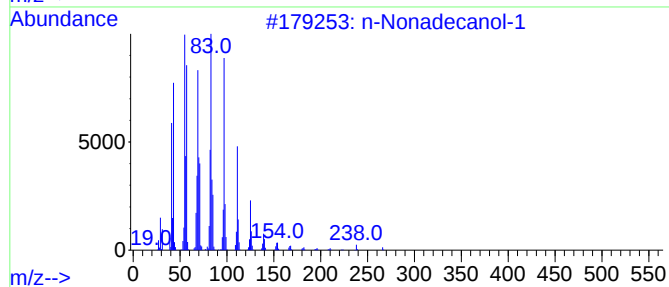
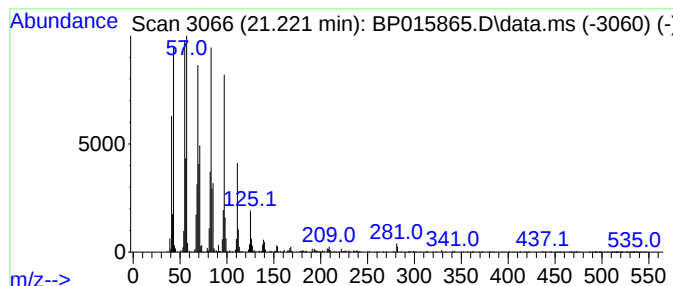
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Nonadecanol-1 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.221	7.86 ng/ul	1230660	Chrysene-d12	21.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Nonadecanol-1	284	C19H40O	001454-84-8	94
2			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
3			n-Tetracosanol-1	354	C24H50O	000506-51-4	94
4			Pentacos-1-ene	350	C25H50	016980-85-1	94
5			Behenic alcohol	326	C22H46O	000661-19-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015865.D
Acq On : 01 Jul 2023 03:02
Operator : MA/JU
Sample : 03161-15
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFX1

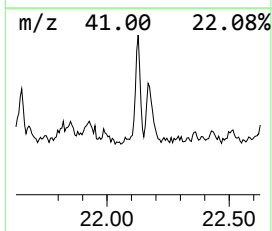
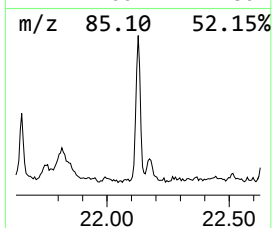
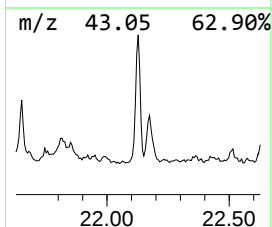
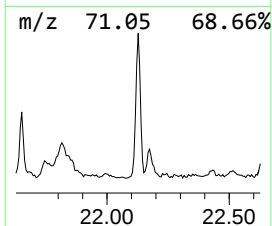
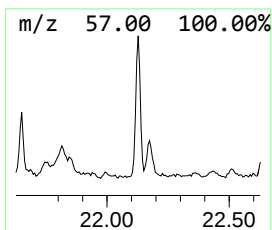
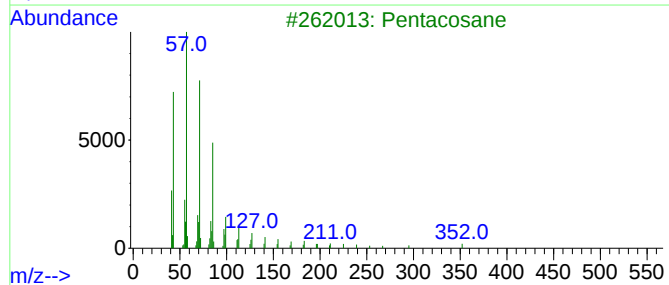
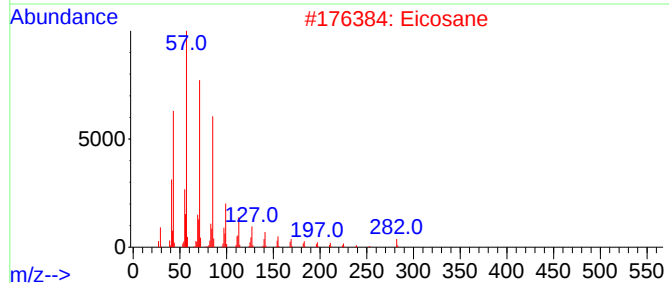
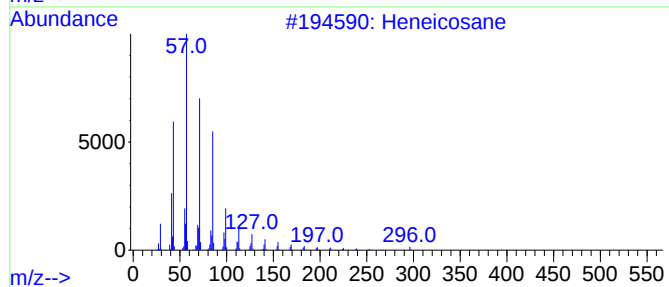
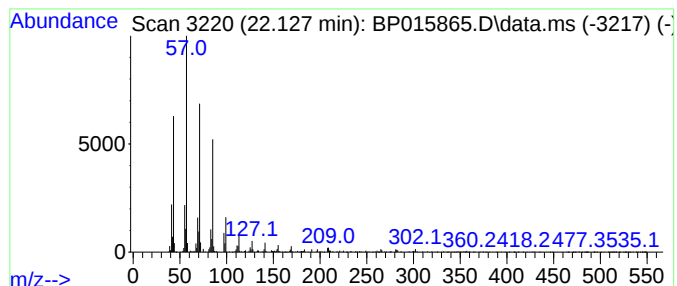
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.127	2.85 ng/ul	446350	Chrysene-d12	21.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	96
2			Eicosane	282	C20H42	000112-95-8	94
3			Pentacosane	352	C25H52	000629-99-2	92
4			Tetracosane	338	C24H50	000646-31-1	87
5			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015865.D
Acq On : 01 Jul 2023 03:02
Operator : MA/JU
Sample : 03161-15
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFX1

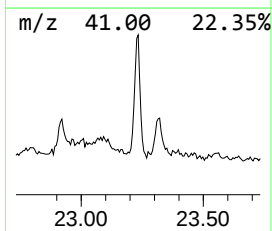
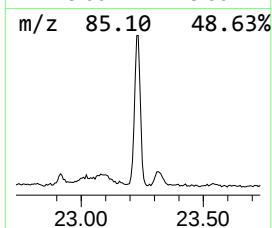
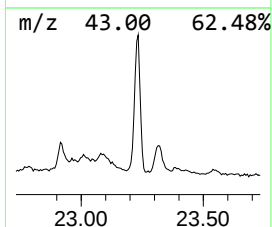
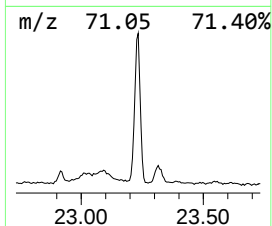
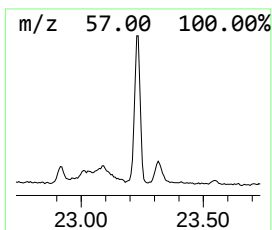
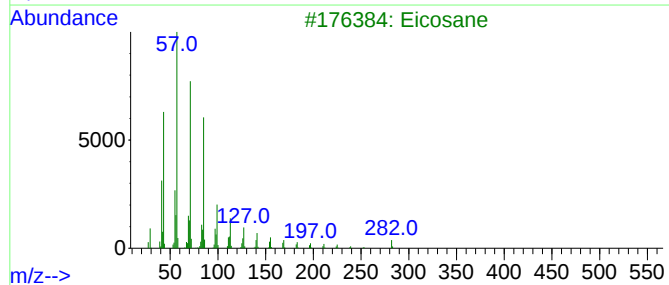
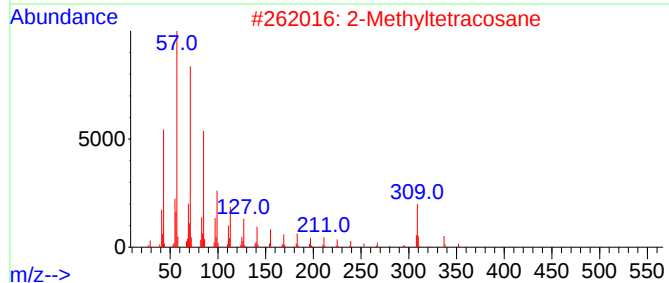
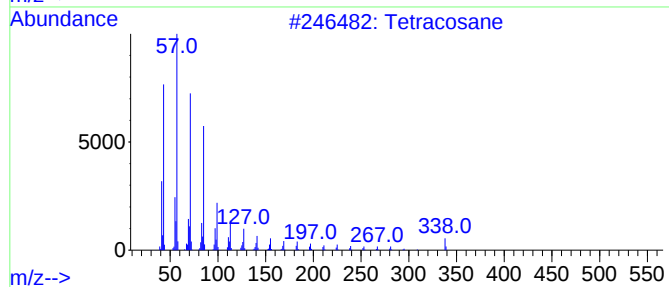
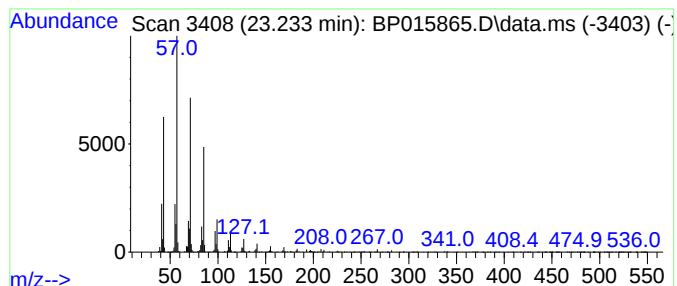
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.233	6.35 ng/ul	1191020	Perylene-d12	24.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	97
2			2-Methyltetracosane	352	C25H52	001560-78-7	96
3			Eicosane	282	C20H42	000112-95-8	93
4			Octacosane	394	C28H58	000630-02-4	91
5			Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015865.D
Acq On : 01 Jul 2023 03:02
Operator : MA/JU
Sample : 03161-15
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFX1

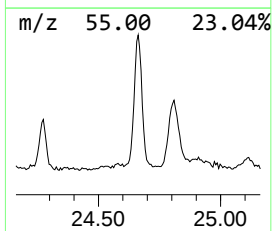
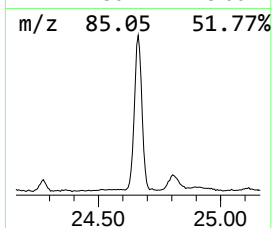
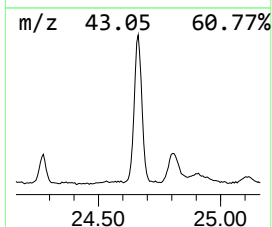
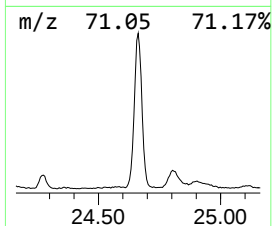
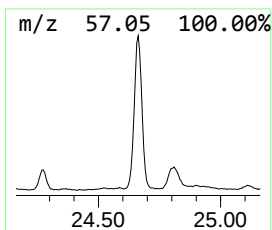
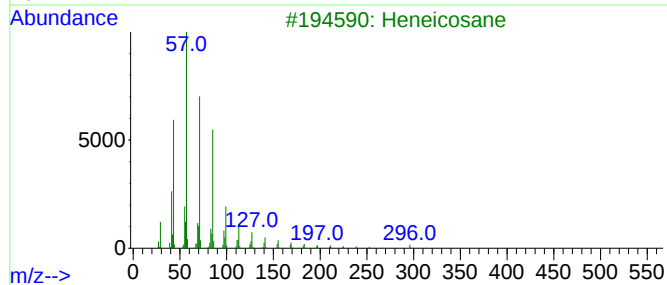
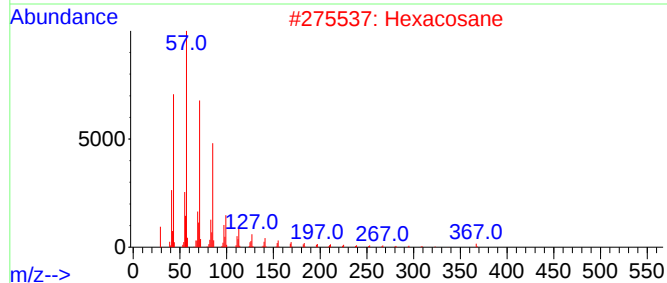
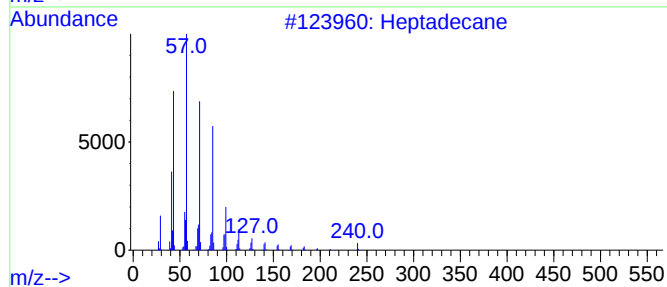
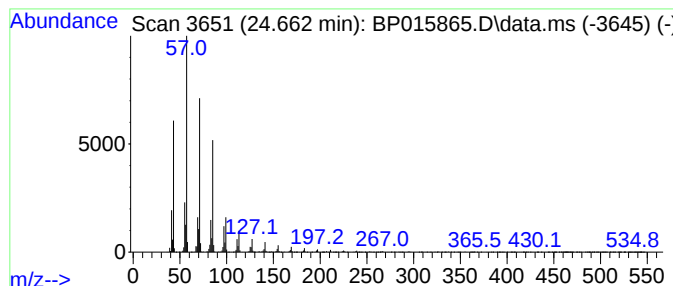
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.662	16.31 ng/ul	3056670	Perylene-d12	24.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	93
2			Hexacosane	366	C26H54	000630-01-3	91
3			Heneicosane	296	C21H44	000629-94-7	91
4			Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91
5			Octacosane	394	C28H58	000630-02-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015865.D
Acq On : 01 Jul 2023 03:02
Operator : MA/JU
Sample : 03161-15
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFX1

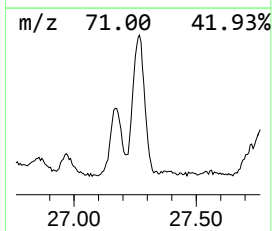
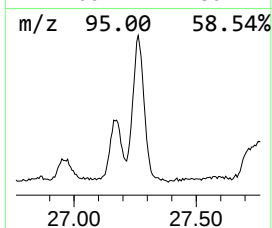
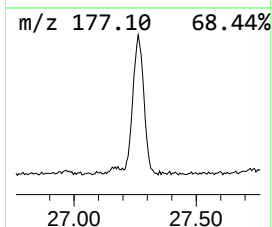
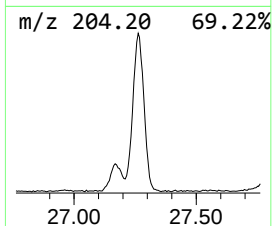
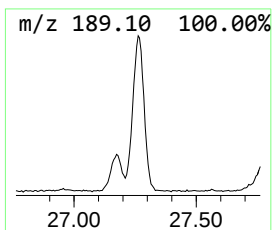
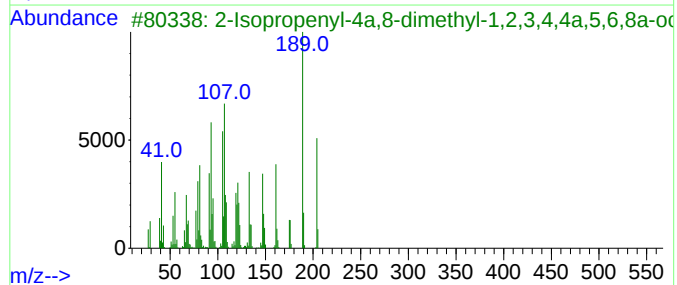
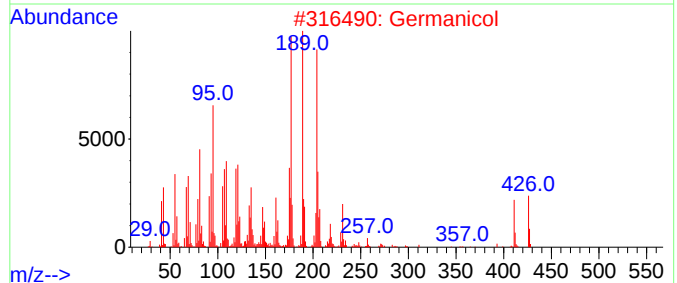
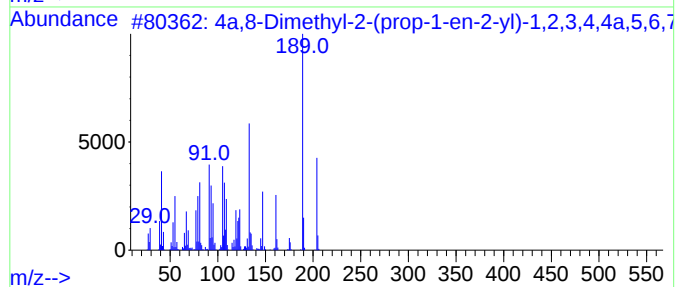
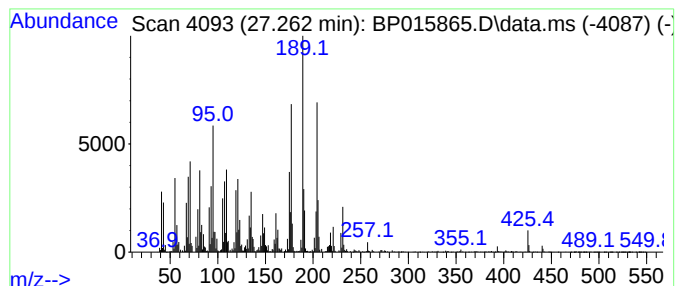
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.262	23.18 ng/ul	4344570	Perylene-d12	24.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	4a,8-Dimethyl-2-(prop-1-en-2-yl)...	204	C15H24	103827-22-1	46		
2	Germanicol	426	C30H50O	000465-02-1	43		
3	2-Isopropenyl-4a,8-dimethyl-1,2,...	204	C15H24	207297-57-2	38		
4	3,4-Dihydrocoumarin, 4,4,7,8-tet...	204	C13H16O2	040614-36-6	30		
5	2,10,10-Trimethyltricyclo[7.1.1...	204	C14H20O	1000210-81-9	30		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015865.D
Acq On : 01 Jul 2023 03:02
Operator : MA/JU
Sample : 03161-15
Misc :
ALS Vial : 31 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFX1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	16.075	3.0	ng/ul	559010	3	14.704	3674160	20.0
cis-10-Nonadece...	18.263	3.1	ng/ul	648812	4	17.463	4166920	20.0
n-Hexadecanoic ...	18.316	10.1	ng/ul	2098260	4	17.463	4166920	20.0
Octadecanoic acid	19.539	3.7	ng/ul	572514	5	21.557	3130020	20.0
Fern-7-en-3.bet...	20.757	5.5	ng/ul	853490	5	21.557	3130020	20.0
n-Nonadecanol-1	21.221	7.9	ng/ul	1230660	5	21.557	3130020	20.0
(DEL) Alkane: S...	22.127	2.9	ng/ul	446350	5	21.557	3130020	20.0
(DEL) Alkane: S...	23.233	6.3	ng/ul	1191020	6	24.098	3749000	20.0
(DEL) Alkane: S...	24.662	16.3	ng/ul	3056670	6	24.098	3749000	20.0
unknown-01	27.262	23.2	ng/ul	4344570	6	24.098	3749000	20.0