

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
 Data File : BP015902.D
 Acq On : 02 Jul 2023 01:35
 Operator : MA/JU
 Sample : 03243-01
 Misc :
 ALS Vial : 70 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCGA5

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: BP015902.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.376	29	32	39	rVB	40448	55264	1.21%	0.089%
2	3.823	106	108	120	rBV	306334	549351	12.00%	0.884%
3	3.923	122	125	134	rVB	59477	93546	2.04%	0.151%
4	4.040	140	145	151	rVB	51237	81135	1.77%	0.131%
5	4.140	160	162	169	rVB2	15724	21441	0.47%	0.035%
6	4.528	226	228	231	rBV4	5135	6779	0.15%	0.011%
7	5.099	320	325	326	rBV2	12997	21700	0.47%	0.035%
8	5.123	326	329	334	rVB2	16148	26964	0.59%	0.043%
9	5.234	343	348	349	rBV4	7319	8873	0.19%	0.014%
10	5.452	383	385	391	rVB7	3796	4989	0.11%	0.008%
11	7.128	665	670	673	rBV7	7540	12722	0.28%	0.020%
12	7.240	682	689	706	rBV	418127	1120908	24.49%	1.804%
13	7.375	706	712	727	rVV	463346	885713	19.35%	1.426%
14	7.587	741	748	773	rVB	593887	1217425	26.60%	1.959%
15	8.046	819	826	842	rBV	831207	1658333	36.24%	2.669%
16	8.693	931	936	938	rBV6	4127	5693	0.12%	0.009%
17	8.793	946	953	974	rBV	466751	1294861	28.30%	2.084%
18	9.246	1021	1030	1051	rBV	536338	1215965	26.57%	1.957%
19	9.581	1084	1087	1091	rVB2	10079	13588	0.30%	0.022%
20	9.975	1145	1154	1174	rBV	365803	1015986	22.20%	1.635%
21	10.546	1242	1251	1274	rBV	416893	1446812	31.62%	2.329%
22	10.881	1300	1308	1324	rVV	1247817	2538863	55.48%	4.086%
23	11.058	1330	1338	1367	rBV	292563	1029068	22.49%	1.656%
24	11.752	1450	1456	1457	rBV6	4205	6210	0.14%	0.010%
25	11.763	1457	1458	1469	rVB6	4430	10857	0.24%	0.017%
26	11.869	1472	1476	1480	rBV7	4977	9575	0.21%	0.015%
27	12.069	1506	1510	1518	rBV9	6499	15832	0.35%	0.025%
28	12.281	1542	1546	1549	rBV6	3136	4924	0.11%	0.008%
29	12.410	1564	1568	1571	rBV6	5287	6886	0.15%	0.011%
30	12.499	1575	1583	1589	rVV9	9883	26207	0.57%	0.042%
31	12.575	1592	1596	1604	rVB9	7201	18514	0.40%	0.030%
32	12.775	1625	1630	1636	rBV7	8148	21281	0.47%	0.034%
33	13.075	1677	1681	1686	rVB7	4024	8170	0.18%	0.013%
34	13.210	1697	1704	1709	rBV10	8126	16549	0.36%	0.027%
35	13.499	1748	1753	1759	rBV2	19395	32009	0.70%	0.052%
36	13.575	1761	1766	1773	rVB2	29860	54471	1.19%	0.088%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
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37	14.016	1834	1841	1845	rBV8	10002	24952	0.55%	0.040%
38	14.110	1850	1857	1872	rBV	1240161	2234235	48.82%	3.596%
39	14.399	1895	1906	1922	rBV	1668798	2791083	60.99%	4.492%
40	14.563	1929	1934	1942	rVB2	28719	46905	1.02%	0.075%
41	14.640	1943	1947	1950	rBV6	4684	7967	0.17%	0.013%
42	14.698	1950	1957	1973	rBV	2068443	3310564	72.34%	5.328%
43	14.922	1990	1995	1999	rBV8	5686	12372	0.27%	0.020%
44	15.004	2002	2009	2025	rBV2	157919	670212	14.65%	1.079%
45	15.110	2025	2027	2034	rVB4	29023	60602	1.32%	0.098%
46	15.316	2060	2062	2064	rBV3	5359	5045	0.11%	0.008%
47	15.463	2085	2087	2091	rVV3	7502	8986	0.20%	0.014%
48	15.516	2093	2096	2103	rVB2	23399	34601	0.76%	0.056%
49	15.698	2121	2127	2142	rBV	1846181	3347797	73.16%	5.388%
50	15.881	2152	2158	2177	rBV2	191739	575493	12.58%	0.926%
51	16.069	2184	2190	2201	rBV	387813	631824	13.81%	1.017%
52	16.345	2235	2237	2240	rVB4	6960	6116	0.13%	0.010%
53	16.398	2240	2246	2252	rBV4	24002	51362	1.12%	0.083%
54	16.551	2268	2272	2278	rBV9	21108	35350	0.77%	0.057%
55	16.840	2319	2321	2327	rVB7	14485	20301	0.44%	0.033%
56	17.175	2372	2378	2381	rBV6	28251	56076	1.23%	0.090%
57	17.334	2401	2405	2412	rBV	94509	158905	3.47%	0.256%
58	17.398	2413	2416	2420	rBV2	56401	68481	1.50%	0.110%
59	17.463	2421	2427	2431	rBV	2410244	3627082	79.26%	5.838%
60	17.504	2431	2434	2439	rVB	284943	409976	8.96%	0.660%
61	17.563	2439	2444	2454	rBV	1932260	3217936	70.32%	5.179%
62	17.910	2499	2503	2509	rVB2	31314	50053	1.09%	0.081%
63	18.063	2526	2529	2532	rBV4	20380	29671	0.65%	0.048%
64	18.204	2548	2553	2557	rBV	80597	123037	2.69%	0.198%
65	18.257	2558	2562	2567	rBV	364664	521514	11.40%	0.839%
66	18.316	2567	2572	2580	rBV	1666835	2215950	48.42%	3.567%
67	19.192	2718	2721	2723	rBV2	44327	59401	1.30%	0.096%
68	19.398	2754	2756	2758	rBV2	64173	73166	1.60%	0.118%
69	19.463	2758	2767	2775	rVB	523094	986336	21.55%	1.588%
70	19.539	2776	2780	2786	rBV	345846	545153	11.91%	0.877%
71	19.792	2818	2823	2836	rVB2	2619306	4576246	100.00%	7.366%
72	20.092	2871	2874	2878	rBV2	125403	146251	3.20%	0.235%
73	20.263	2900	2903	2907	rBV	139710	172556	3.77%	0.278%
74	20.751	2983	2986	2989	rBV4	149408	270230	5.91%	0.435%
75	21.222	3060	3066	3072	rBV3	648599	1117024	24.41%	1.798%
76	21.410	3095	3098	3102	rBV	335789	379199	8.29%	0.610%

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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

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Peak separation: 5

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77	21.551	3117	3122	3127	rBV2	1944077	2959836	64.68%	4.764%
78	23.227	3402	3407	3415	rVB	1016276	1602624	35.02%	2.579%
79	23.927	3522	3526	3534	rVB2	1140426	2146004	46.89%	3.454%
80	24.092	3549	3554	3570	rVB	1324625	3437133	75.11%	5.532%
81	24.545	3626	3631	3642	rVB	1458180	3171531	69.30%	5.105%
82	24.657	3645	3650	3662	rVB	740156	1574954	34.42%	2.535%

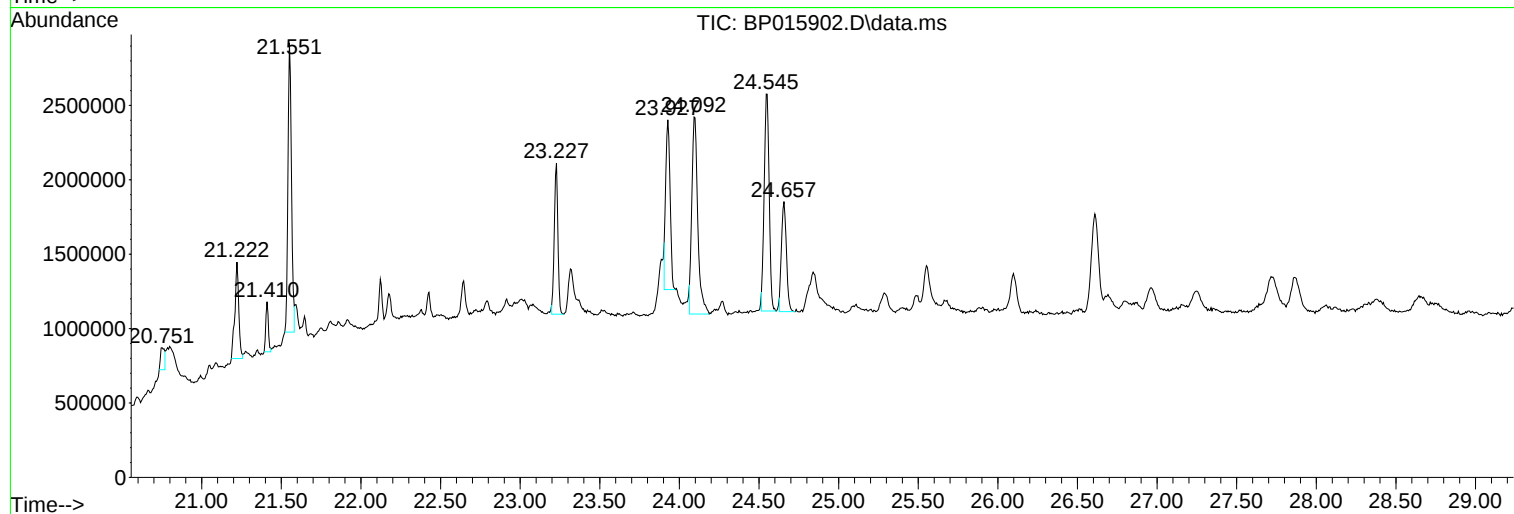
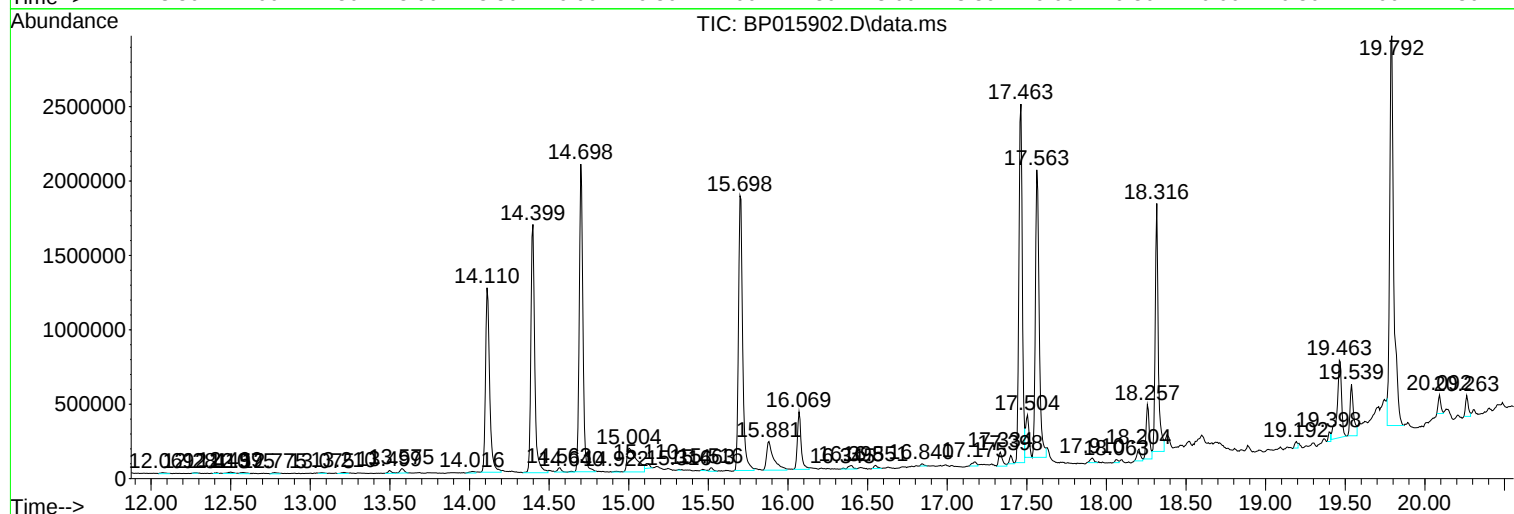
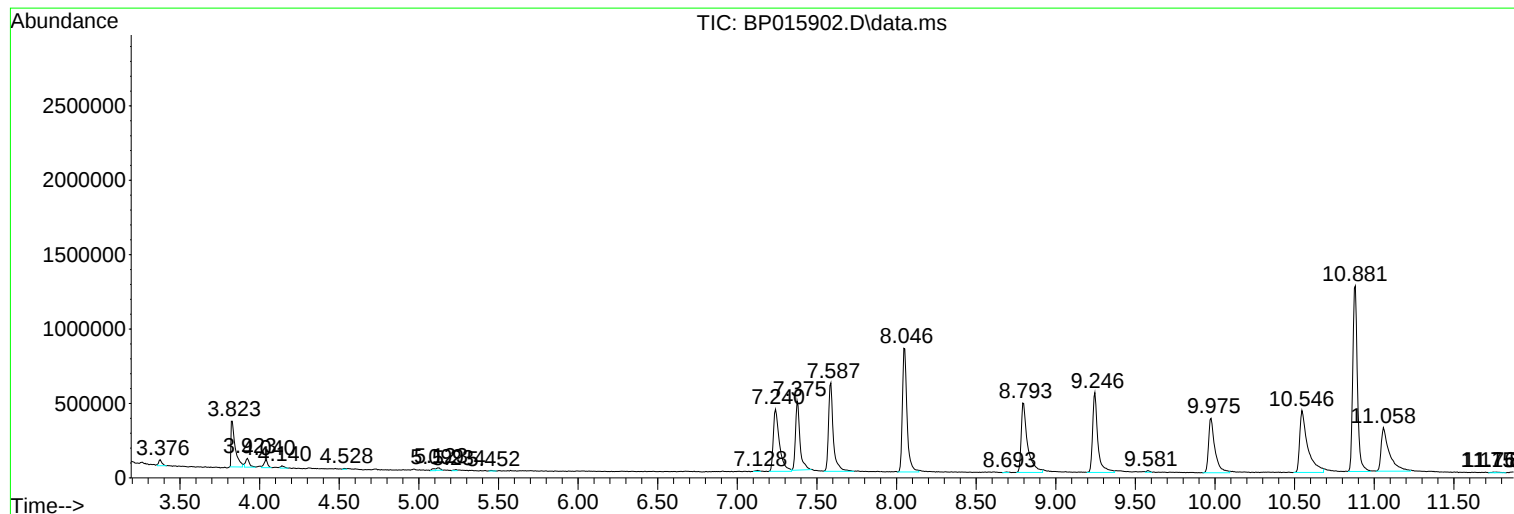
Sum of corrected areas: 62129556

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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



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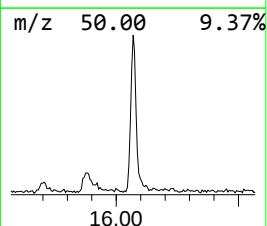
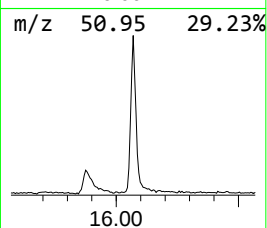
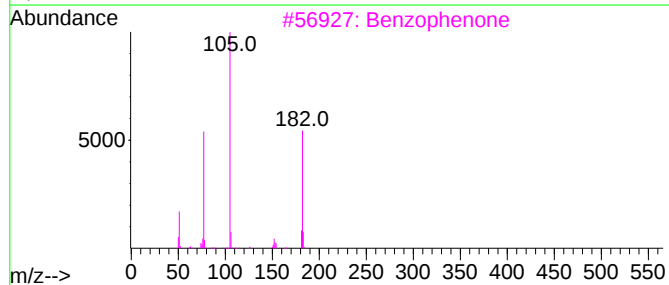
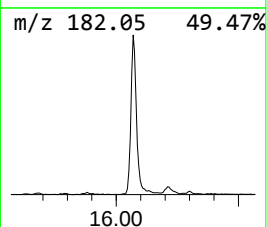
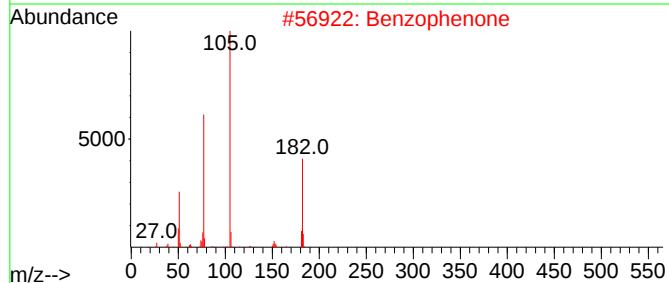
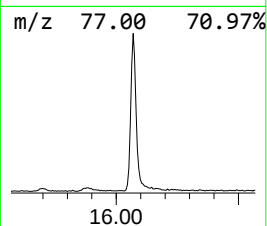
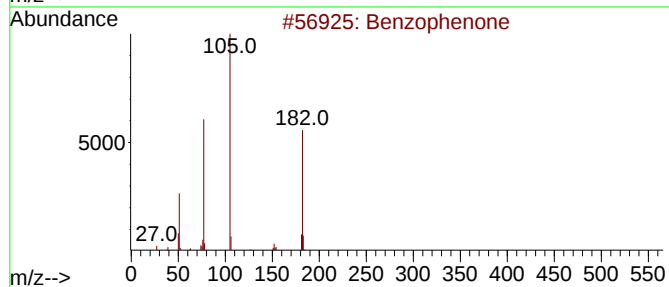
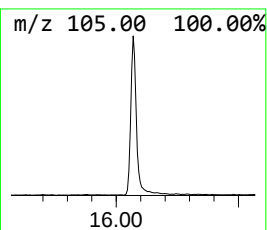
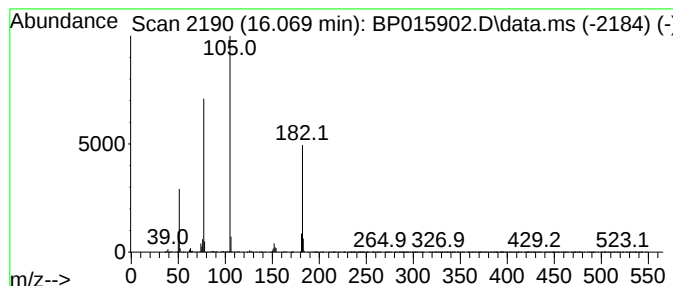
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.069	3.82 ng/ul	631824	Acenaphthene-d10	14.698

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	95
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	86



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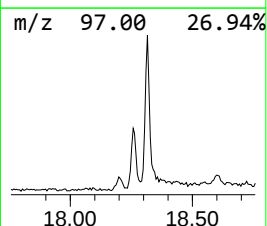
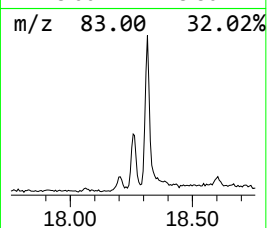
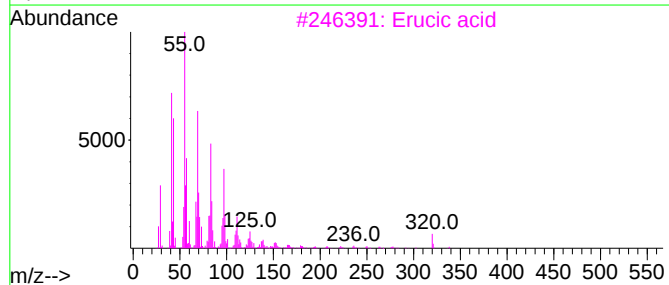
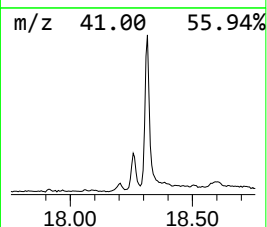
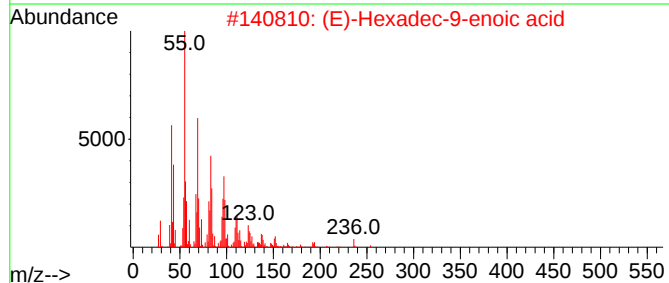
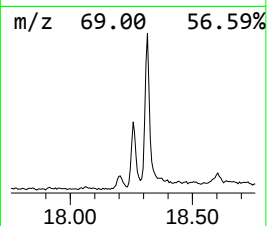
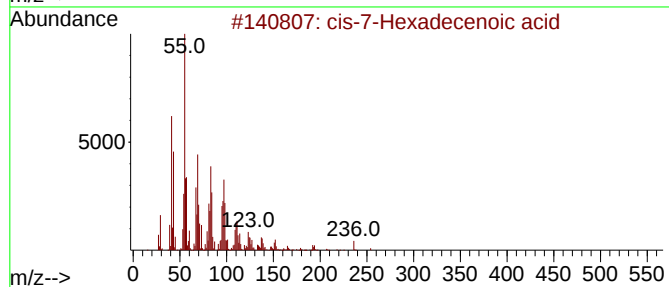
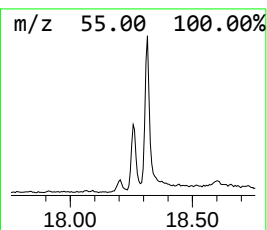
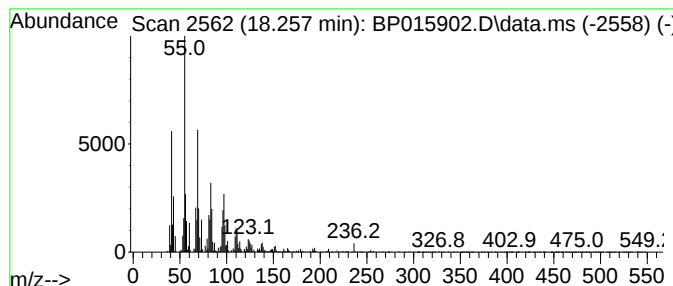
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 cis-7-Hexadecenoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.257	2.88 ng/ul	521514	Phenanthrene-d10	17.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cis-7-Hexadecenoic acid	254	C16H30O2	002416-19-5	98
2			(E)-Hexadec-9-enoic acid	254	C16H30O2	010030-73-6	98
3			Erucic acid	338	C22H42O2	000112-86-7	97
4			Palmitoleic acid	254	C16H30O2	000373-49-9	96
5			Palmitoleic acid	254	C16H30O2	000373-49-9	92



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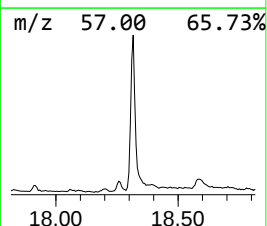
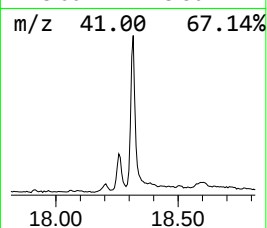
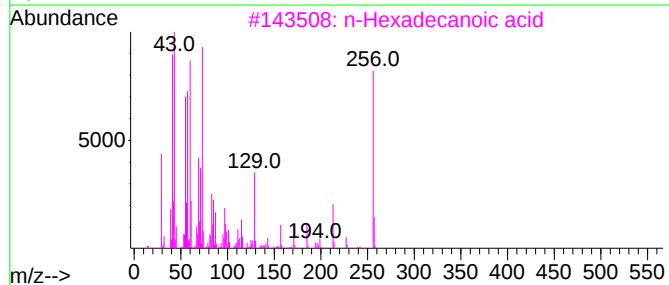
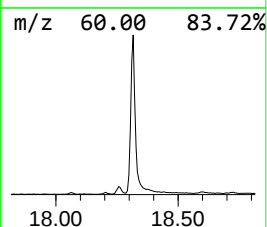
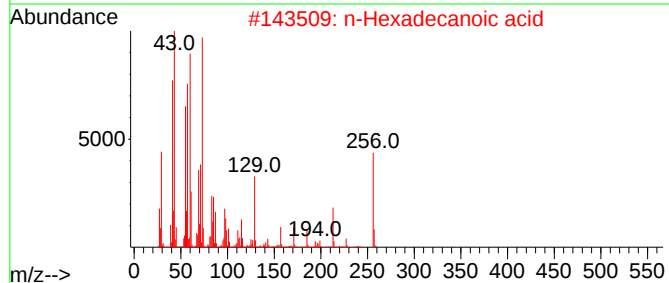
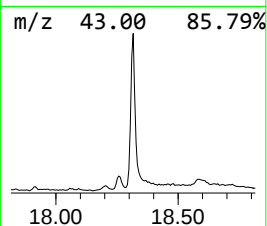
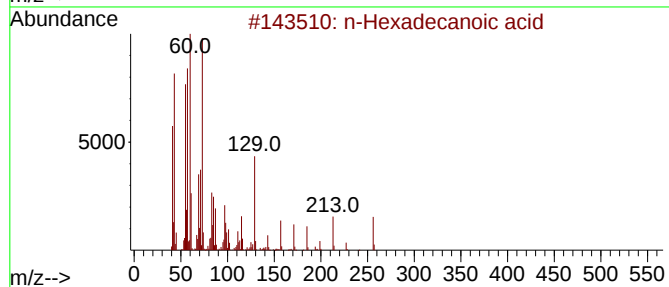
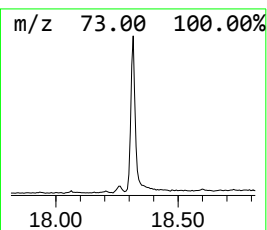
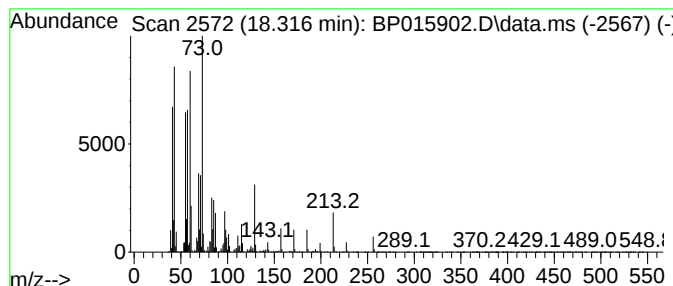
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.316	12.22 ng/ul	2215950	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	99
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	94



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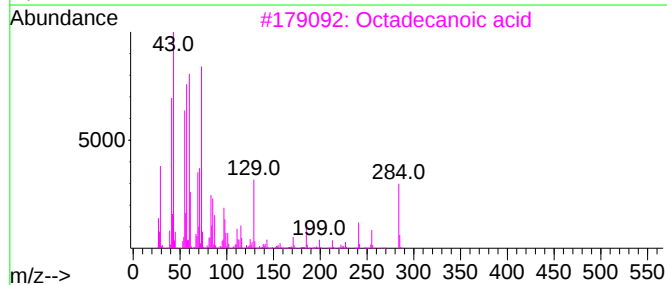
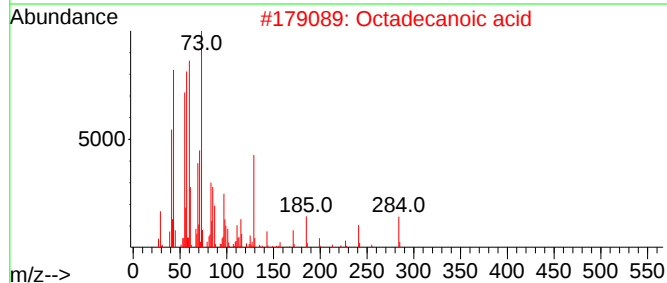
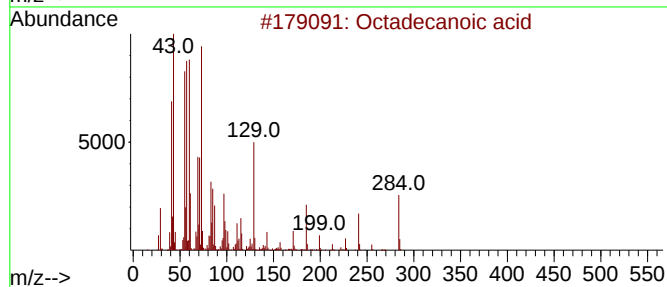
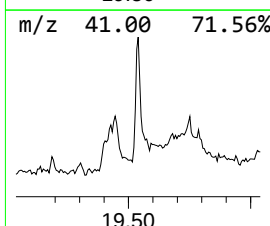
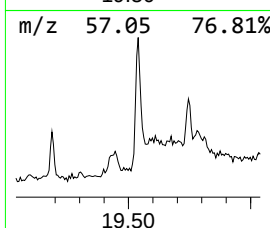
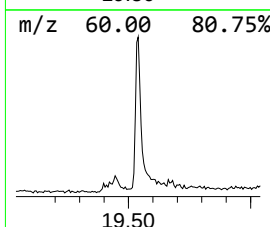
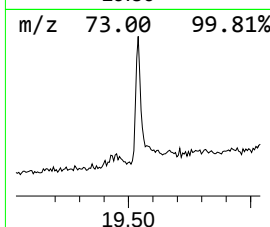
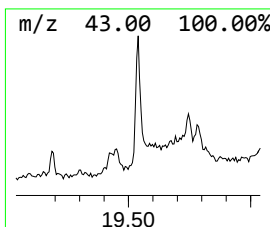
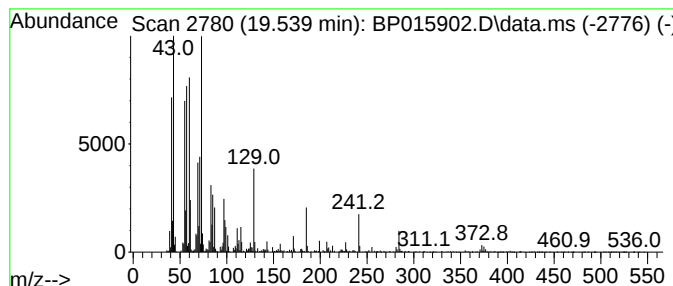
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.539	3.68 ng/ul	545153	Chrysene-d12	21.551

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	95
4			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	94
5			Octadecanoic acid	284	C18H36O2	000057-11-4	92



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015902.D
Acq On : 02 Jul 2023 01:35
Operator : MA/JU
Sample : 03243-01
Misc :
ALS Vial : 70 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGA5

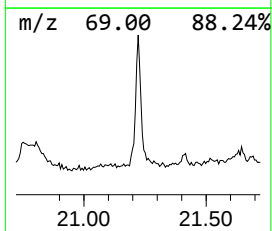
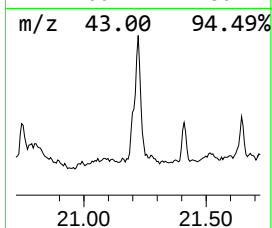
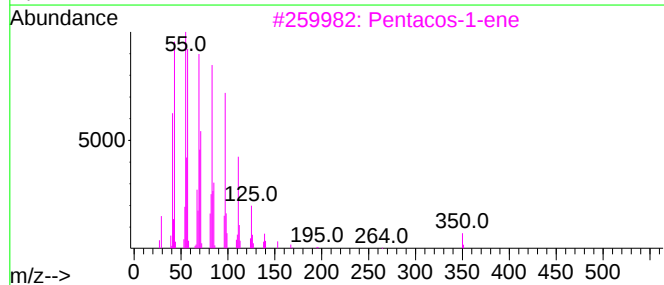
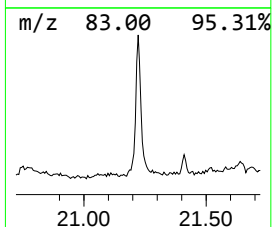
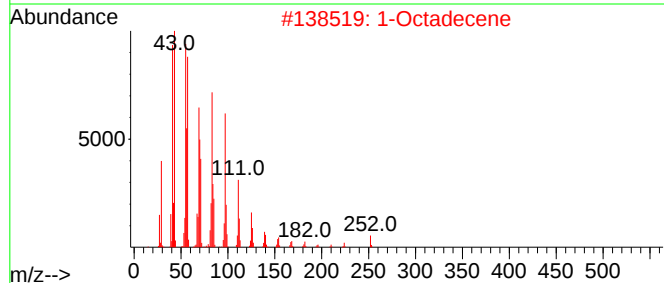
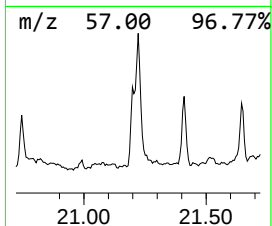
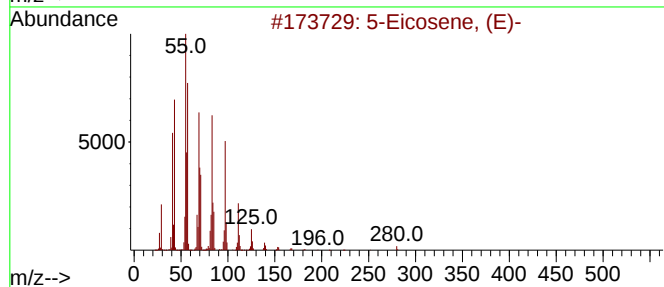
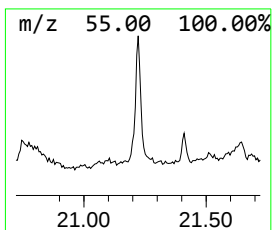
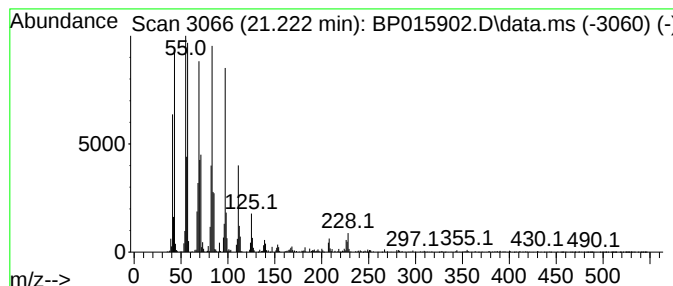
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.222	7.55 ng/ul	1117020	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-			280	C20H40	074685-30-6	99
2	1-Octadecene			252	C18H36	000112-88-9	95
3	Pentacos-1-ene			350	C25H50	016980-85-1	94
4	1-Octadecanol			270	C18H38O	000112-92-5	94
5	Behenic alcohol			326	C22H46O	000661-19-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015902.D
Acq On : 02 Jul 2023 01:35
Operator : MA/JU
Sample : 03243-01
Misc :
ALS Vial : 70 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGA5

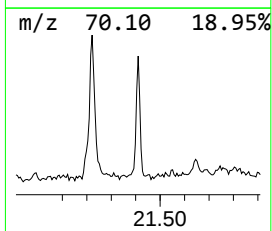
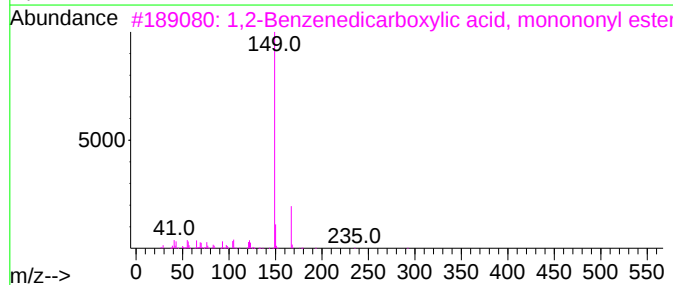
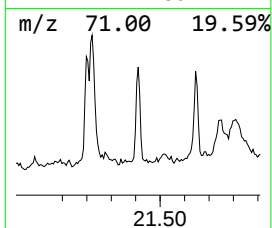
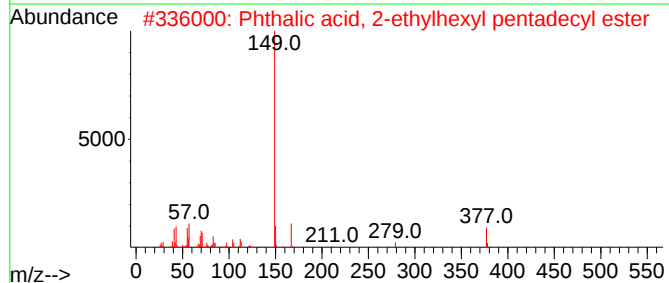
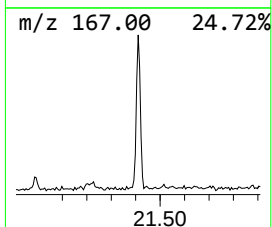
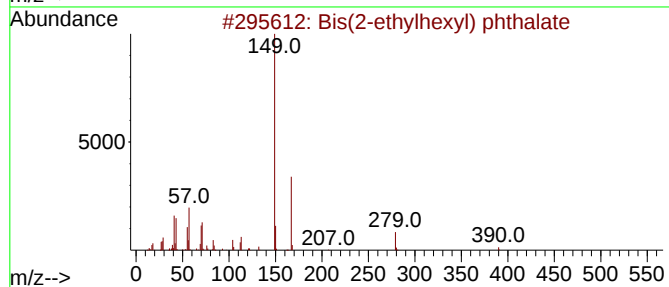
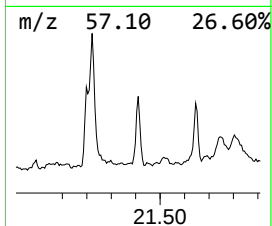
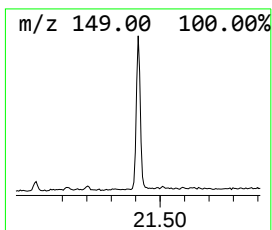
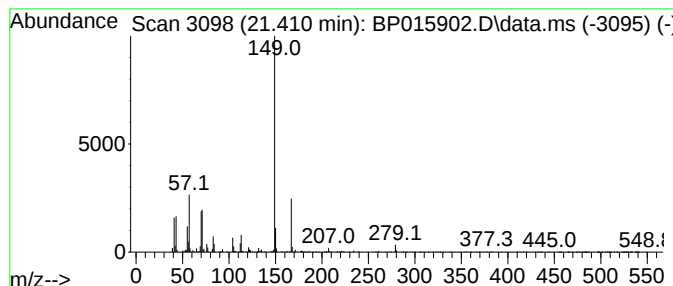
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Bis(2-ethylhexyl) phthalate Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.410	2.56 ng/ul	379199	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	86
2			Phthalic acid, 2-ethylhexyl pent...	488	C31H52O4	1000308-98-2	64
3			1,2-Benzenedicarboxylic acid, mo...	292	C17H24O4	024539-59-1	64
4			Phthalic acid, di(hex-3-yl) ester	334	C20H30O4	1000356-96-7	59
5			Phthalic acid, di(2-methylbutyl)...	306	C18H26O4	1000322-82-5	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015902.D
Acq On : 02 Jul 2023 01:35
Operator : MA/JU
Sample : 03243-01
Misc :
ALS Vial : 70 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGA5

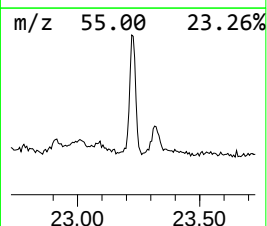
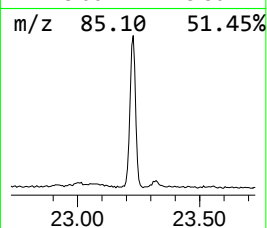
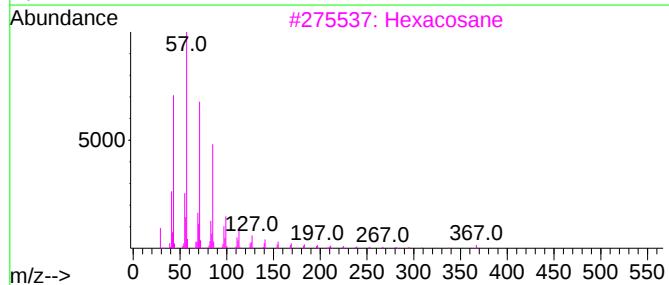
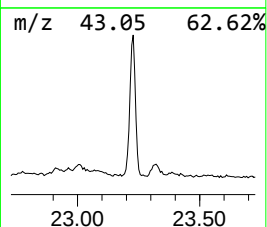
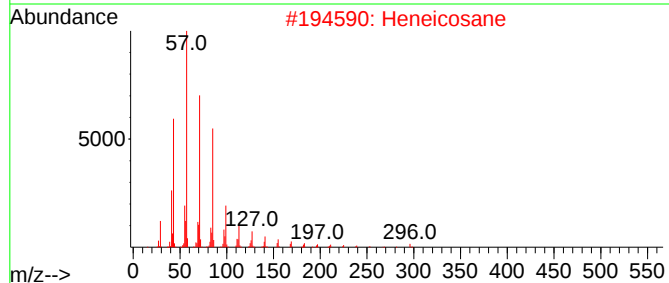
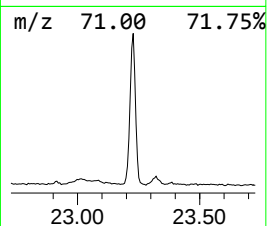
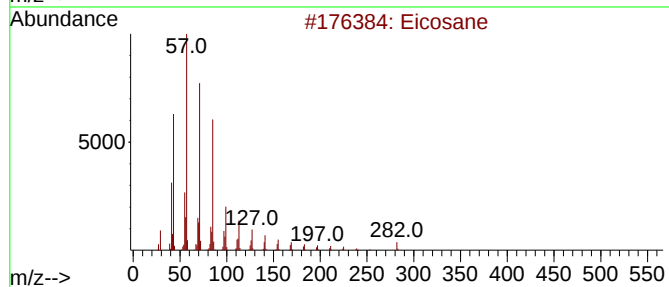
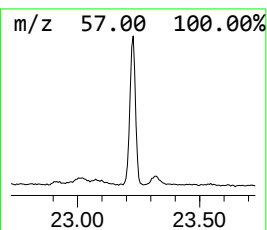
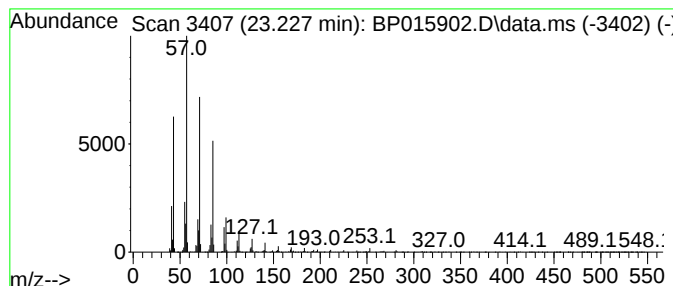
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.227	9.33 ng/ul	1602620	Perylene-d12	24.092

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane	282	C20H42	000112-95-8	96
2		Heneicosane	296	C21H44	000629-94-7	91
3		Hexacosane	366	C26H54	000630-01-3	91
4		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
5		Hexatriacontane	507	C36H74	000630-06-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015902.D
Acq On : 02 Jul 2023 01:35
Operator : MA/JU
Sample : 03243-01
Misc :
ALS Vial : 70 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGA5

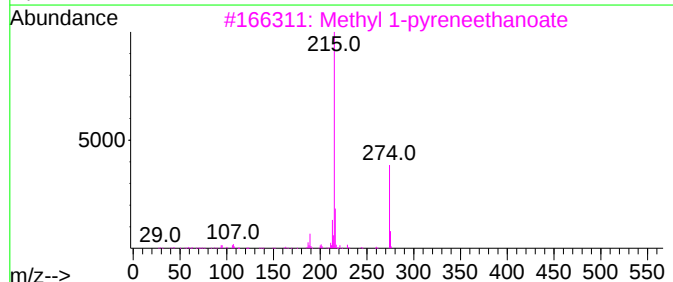
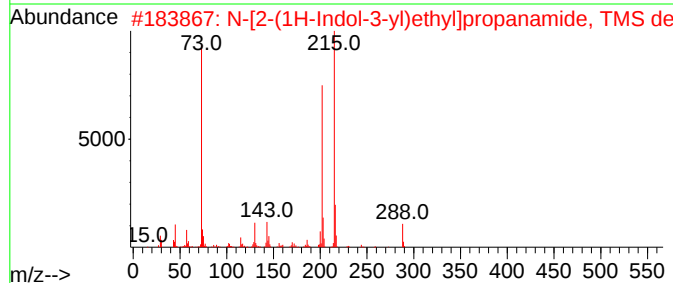
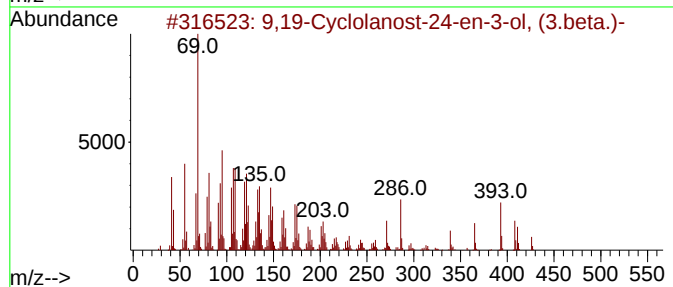
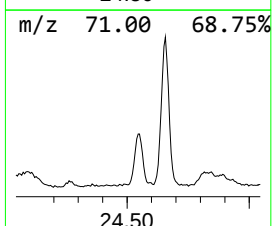
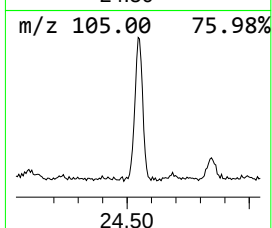
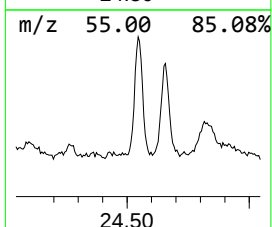
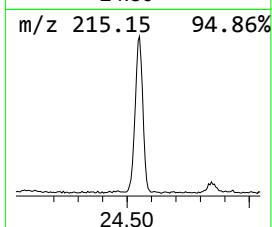
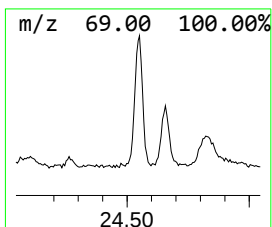
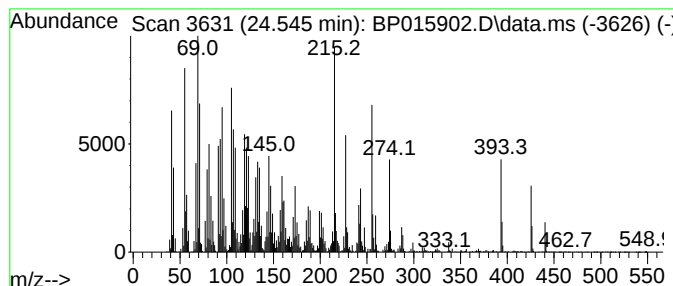
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.545	18.45 ng/ul	3171530	Perylene-d12	24.092

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,19-Cyclolanost-24-en-3-ol, (3...	426	C30H50O	000469-38-5	42		
2	N-[2-(1H-Indol-3-yl)ethyl]propan...	288	C16H24N2OSi	1000493-19-0	25		
3	Methyl 1-pyreneethanoate	274	C19H14O2	073654-19-0	20		
4	Chlorflurecol methyl ester	274	C15H11ClO3	002536-31-4	18		
5	6H-Benzo[d]pyrimido[5,4-b]thiopy...	215	C11H9N3S	1010267-44-6	15		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015902.D
Acq On : 02 Jul 2023 01:35
Operator : MA/JU
Sample : 03243-01
Misc :
ALS Vial : 70 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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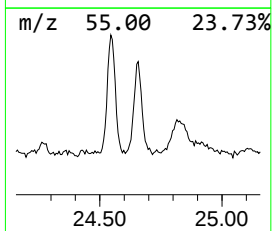
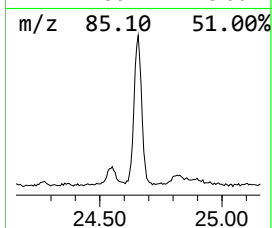
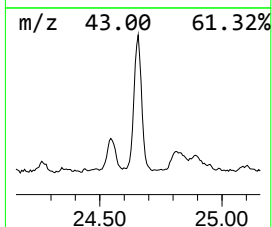
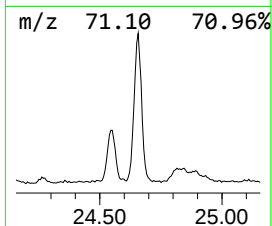
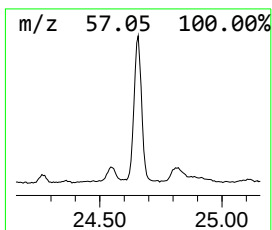
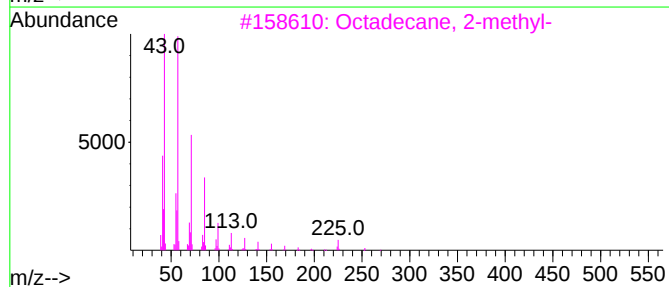
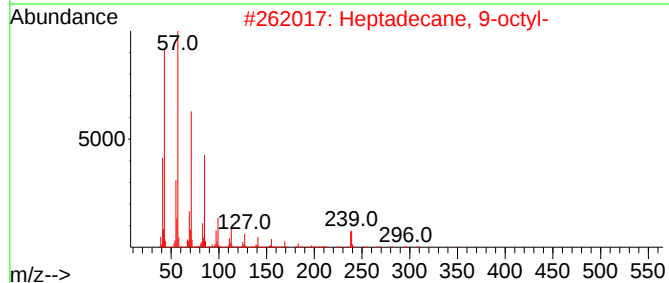
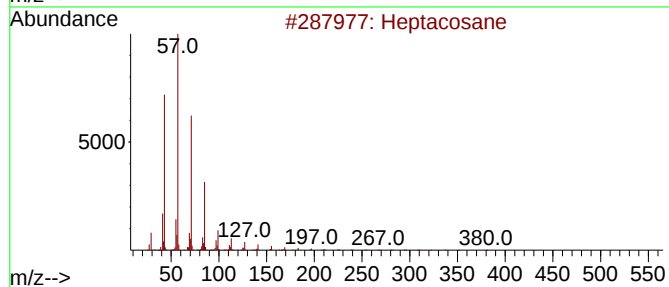
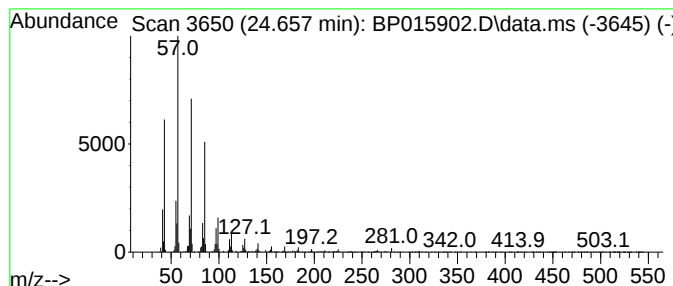
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.657	9.16 ng/ul	1574950	Perylene-d12	24.092

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptacosane	380	C27H56	000593-49-7	94
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	94
3		Octadecane, 2-methyl-	268	C19H40	001560-88-9	93
4		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	91
5		Octacosane	394	C28H58	000630-02-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP063023\
Data File : BP015902.D
Acq On : 02 Jul 2023 01:35
Operator : MA/JU
Sample : 03243-01
Misc :
ALS Vial : 70 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCGA5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.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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cis-7-Hexadecen...	18.257	2.9	ng/u1	521514	4	17.463	3627080	20.0
n-Hexadecanoic ...	18.316	12.2	ng/u1	2215950	4	17.463	3627080	20.0
Octadecanoic acid	19.539	3.7	ng/u1	545153	5	21.551	2959840	20.0
(DEL) Alkane: S...	21.222	7.5	ng/u1	1117020	5	21.551	2959840	20.0
Bis(2-ethylhexy...	21.410	2.6	ng/u1	379199	5	21.551	2959840	20.0
(DEL) Alkane: S...	23.227	9.3	ng/u1	1602620	6	24.092	3437130	20.0
unknown-01	24.545	18.4	ng/u1	3171530	6	24.092	3437130	20.0
(DEL) Alkane: S...	24.657	9.2	ng/u1	1574950	6	24.092	3437130	20.0