

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
 Data File : BM043761.D
 Acq On : 05 Jan 2024 15:37
 Operator : MA/JU
 Sample : 05919-15ME
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 GCF47ME

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_M\Methods\SFAM-EPA-BM121523.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM043761.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.798	101	104	117	rBV	142852	257770	0.26%	0.138%
2	7.128	662	670	682	rBV	672764	1147925	1.18%	0.616%
3	7.292	692	698	707	rBV	546942	866713	0.89%	0.465%
4	7.481	723	730	738	rBV	686679	1121796	1.15%	0.602%
5	7.951	803	810	817	rBV	941576	1507093	1.54%	0.809%
6	8.675	926	933	945	rBV	772423	1352346	1.38%	0.726%
7	9.128	1003	1010	1022	rBV	596784	1050358	1.08%	0.564%
8	9.851	1122	1133	1141	rBV	509612	903572	0.92%	0.485%
9	10.392	1215	1225	1246	rBV	700285	1376495	1.41%	0.738%
10	10.763	1279	1288	1301	rBV	1179356	2048441	2.10%	1.099%
11	10.922	1308	1315	1332	rBV	288148	681860	0.70%	0.366%
12	13.121	1684	1689	1694	rVB	148933	201614	0.21%	0.108%
13	13.392	1731	1735	1746	rBV6	51630	132804	0.14%	0.071%
14	14.015	1836	1841	1846	rBV	1318256	1895742	1.94%	1.017%
15	14.068	1846	1850	1854	rVB	387262	525003	0.54%	0.282%
16	14.292	1882	1888	1894	rBV	1675052	2583684	2.64%	1.386%
17	14.421	1906	1910	1915	rVB3	127125	222310	0.23%	0.119%
18	14.480	1915	1920	1926	rBV	2587572	3610529	3.70%	1.937%
19	14.551	1929	1932	1934	rVV2	121104	166129	0.17%	0.089%
20	14.598	1934	1940	1946	rVB	1752928	2765416	2.83%	1.484%
21	14.833	1975	1980	1986	rBV2	611562	981950	1.01%	0.527%
22	14.986	2003	2006	2010	rBV	162271	220413	0.23%	0.118%
23	15.104	2022	2026	2030	rBV2	284971	391019	0.40%	0.210%
24	15.145	2030	2033	2037	rVV4	130025	177009	0.18%	0.095%
25	15.227	2038	2047	2054	rVB	1032457	1830272	1.87%	0.982%
26	15.421	2075	2080	2087	rBV2	830113	1371374	1.40%	0.736%
27	15.592	2104	2109	2115	rBV	2030723	3008869	3.08%	1.614%
28	15.704	2122	2128	2141	rVB	6743560	10041423	10.28%	5.387%
29	15.839	2146	2151	2158	rVV	1351791	2220968	2.27%	1.192%
30	16.321	2227	2233	2246	rVB	3766298	6436632	6.59%	3.453%
31	16.562	2269	2274	2279	rBV2	851691	1367479	1.40%	0.734%
32	17.021	2343	2352	2357	rBV	44464869	97685134	100.00%	52.407%
33	17.151	2369	2374	2379	rVB	4115967	6066318	6.21%	3.255%
34	17.351	2403	2408	2415	rBV	2333198	3988733	4.08%	2.140%
35	17.451	2421	2425	2433	rVB	2381072	3415326	3.50%	1.832%
36	17.756	2473	2477	2483	rVB	1508977	2276408	2.33%	1.221%

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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_M\Methods\SFAM-EPA-BM121523.MA.M
Title : SVOA CALIBRATION

37	18.245	2557	2560	2564	rBV	1047979	1430341	1.46%	0.767%
38	19.739	2810	2814	2818	rBV	2792515	3847798	3.94%	2.064%
39	19.839	2829	2831	2841	rVB	942901	1416619	1.45%	0.760%
40	21.239	3066	3069	3077	rVB3	682930	863165	0.88%	0.463%
41	21.533	3115	3119	3128	rVB	2706658	3738236	3.83%	2.006%
42	23.791	3498	3503	3510	rVB2	1967795	3757452	3.85%	2.016%
43	23.950	3525	3530	3538	rVB	1750485	3667970	3.75%	1.968%
44	25.668	3818	3822	3833	rVB2	807388	1778292	1.82%	0.954%

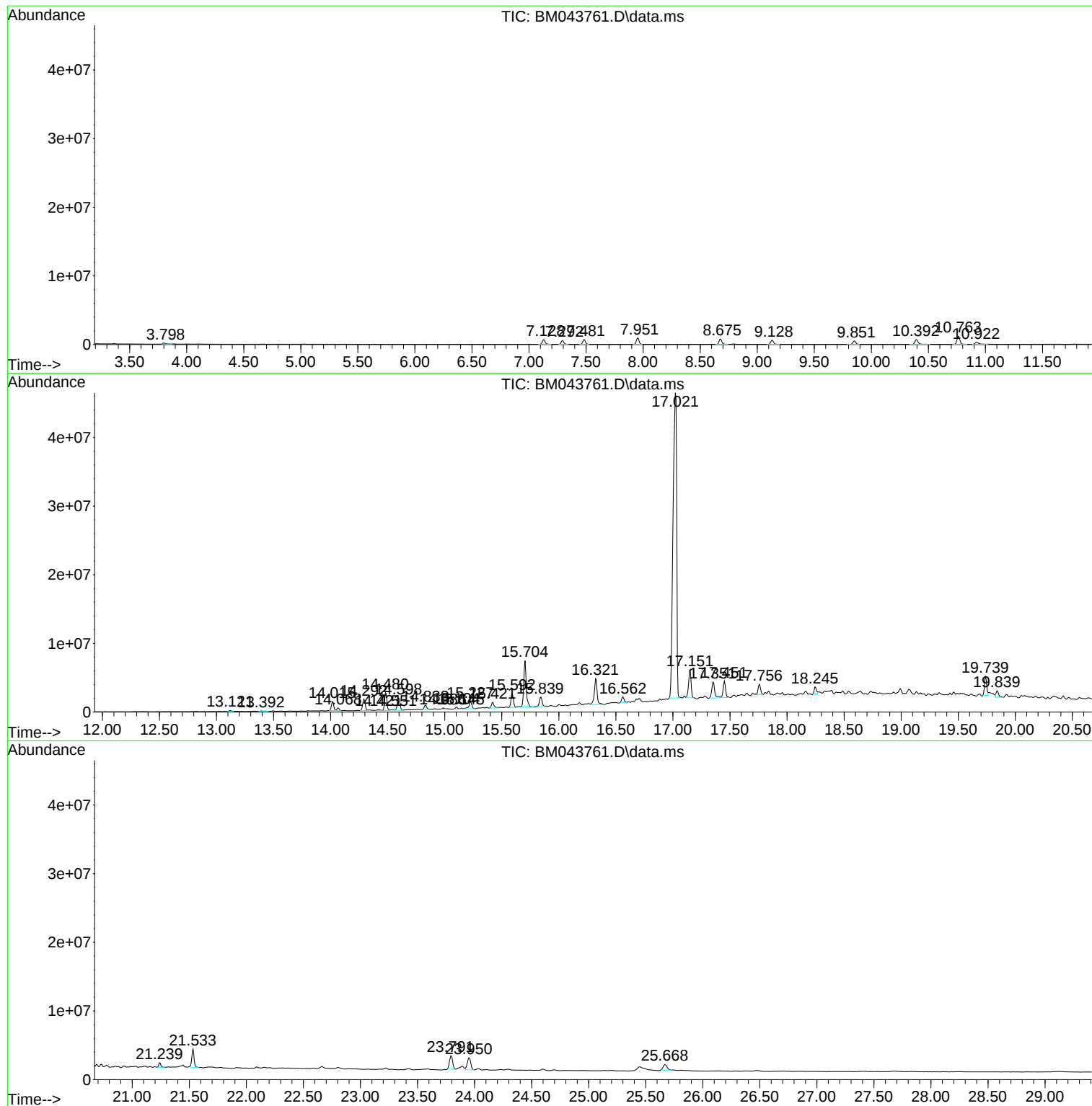
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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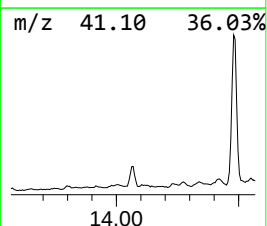
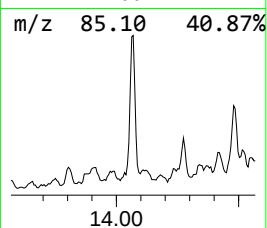
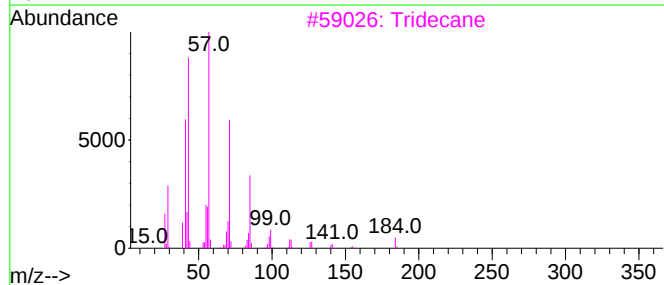
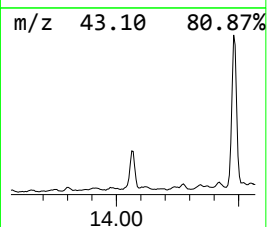
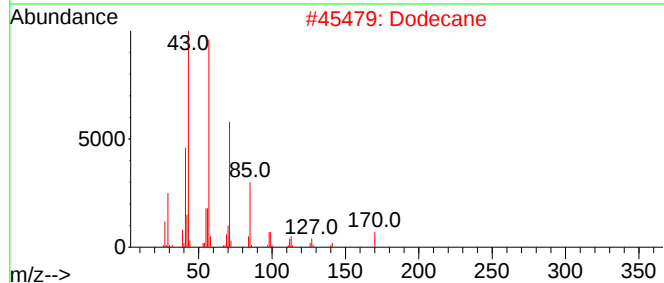
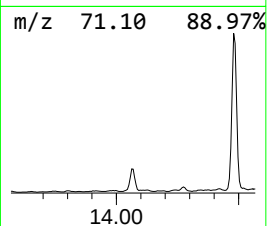
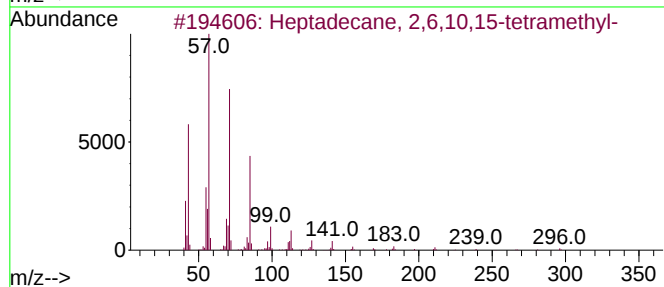
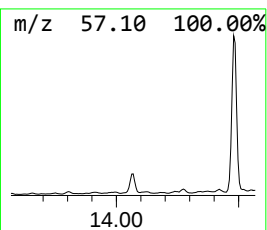
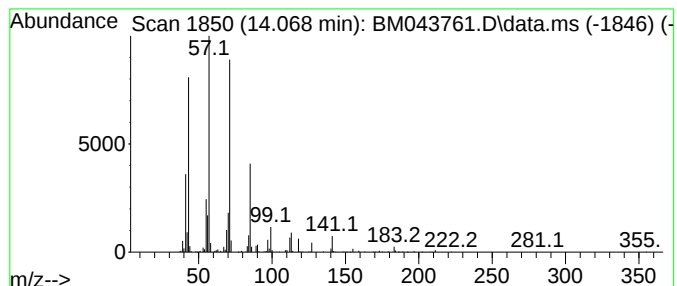
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.068	3.80 ng/ul	525003	Acenaphthene-d10	14.598

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
2		Dodecane	170	C12H26	000112-40-3	90
3		Tridecane	184	C13H28	000629-50-5	90
4		Decane, 5-propyl-	184	C13H28	017312-62-8	87
5		Tetradecane	198	C14H30	000629-59-4	86



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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
GCF47ME

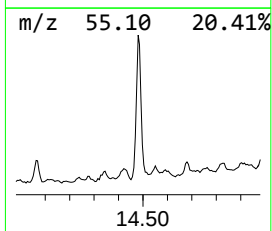
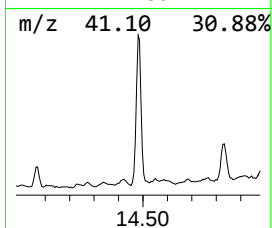
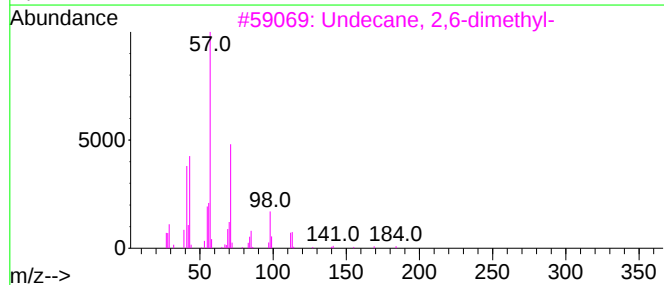
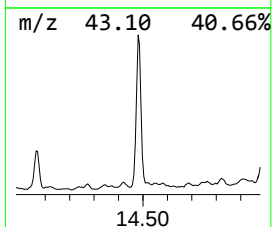
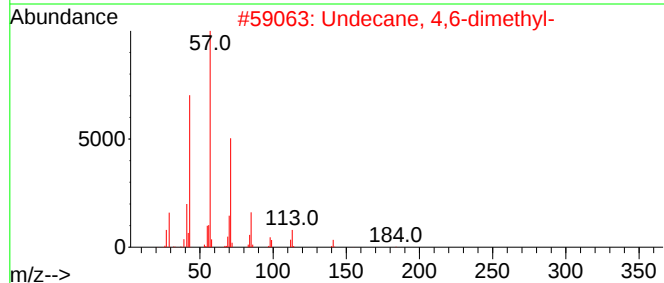
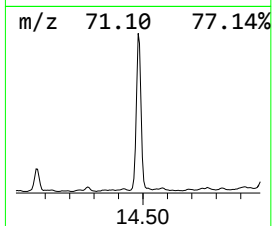
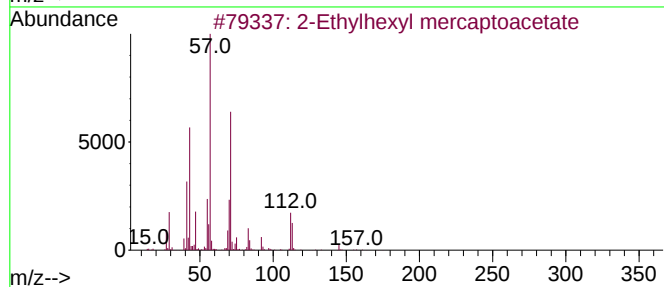
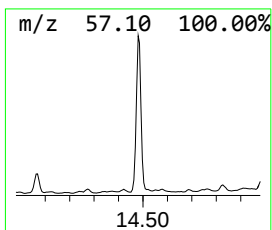
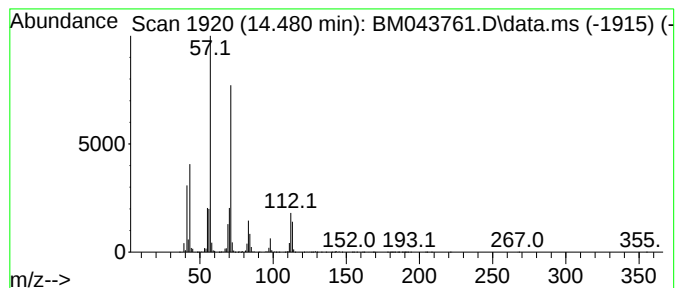
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2-Ethylhexyl mercaptoacetate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.480	26.11 ng/ul	3610530	Acenaphthene-d10	14.598

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Ethylhexyl mercaptoacetate	204	C10H20O2S	007659-86-1	72
2			Undecane, 4,6-dimethyl-	184	C13H28	017312-82-2	72
3			Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	72
4			1-Decene, 3,4-dimethyl-	168	C12H24	050871-03-9	72
5			Carbonic acid, neopentyl 2-ethyl...	244	C14H28O3	1000357-85-7	64



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Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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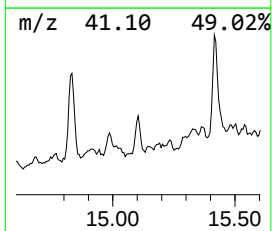
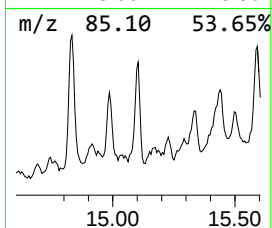
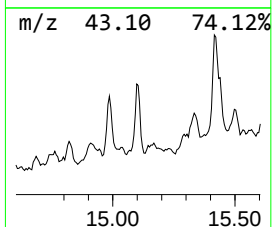
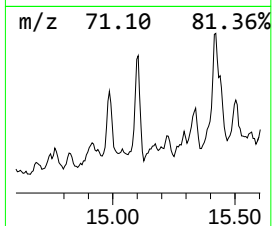
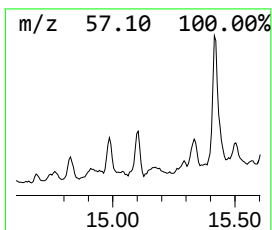
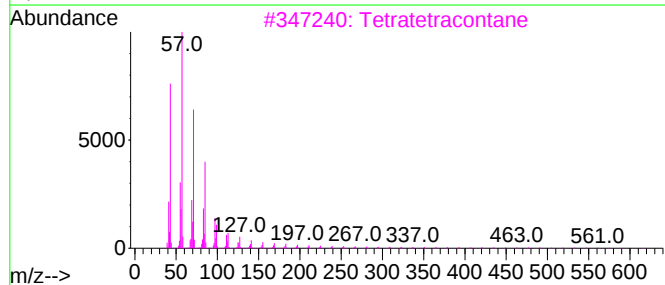
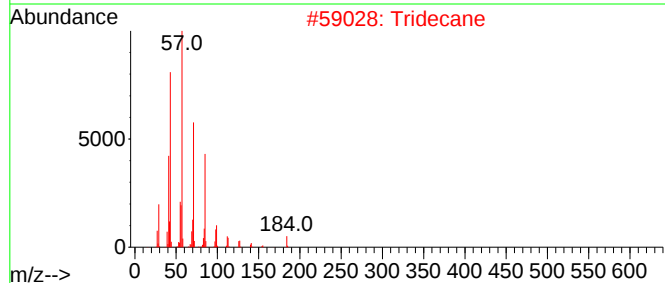
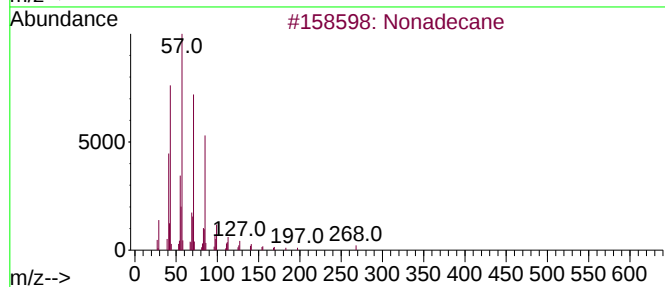
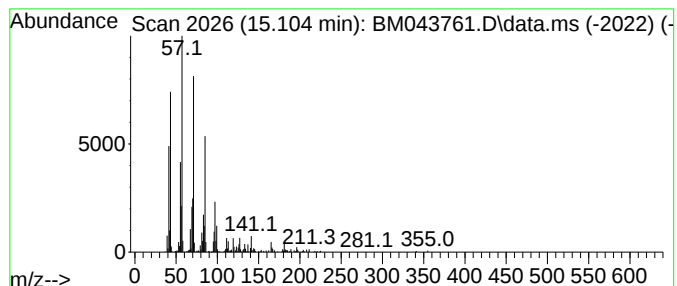
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.104	2.83 ng/ul	391019	Acenaphthene-d10	14.598

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	81
2			Tridecane	184	C13H28	000629-50-5	81
3			Tetratetracontane	619	C44H90	007098-22-8	80
4			Hexacosane	366	C26H54	000630-01-3	72
5			Heptacosane	380	C27H56	000593-49-7	72



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Instrument :
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ClientSampleId :
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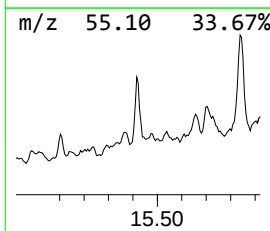
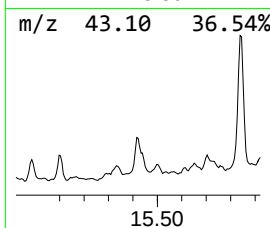
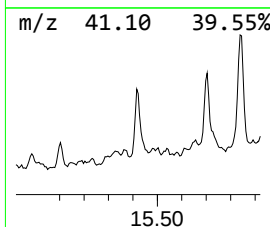
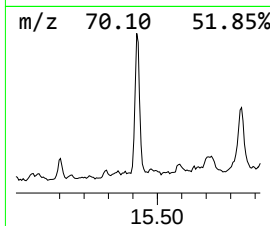
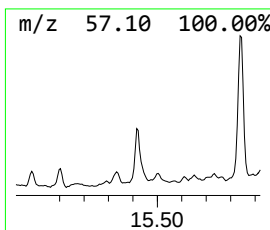
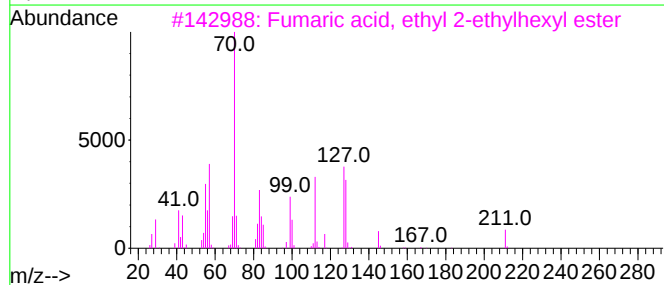
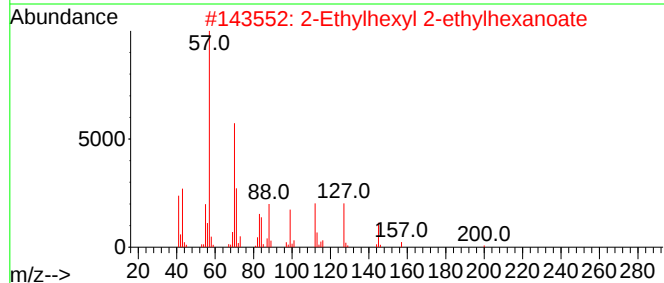
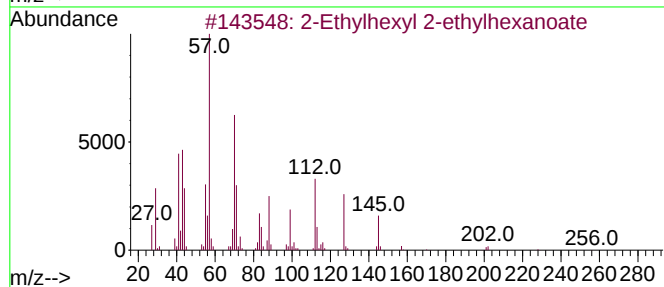
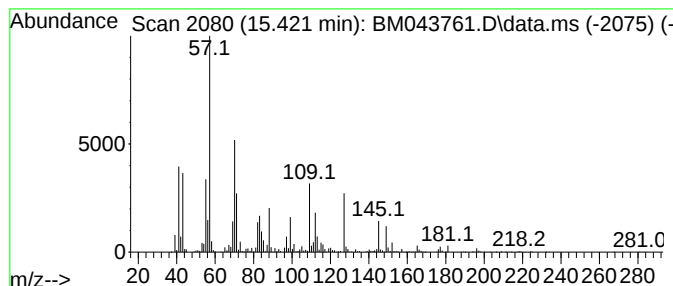
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 2-Ethylhexyl 2-ethylhexanoate Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.421	9.92 ng/ul	1371370	Acenaphthene-d10	14.598

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Ethylhexyl 2-ethylhexanoate	256	C16H32O2	007425-14-1	53
2			2-Ethylhexyl 2-ethylhexanoate	256	C16H32O2	007425-14-1	52
3			Fumaric acid, ethyl 2-ethylhexyl...	256	C14H24O4	1000339-13-5	46
4			Octanoic acid, 2-ethylhexyl ester	256	C16H32O2	063321-70-0	43
5			Acetic acid, 2-ethylhexyl ester	172	C10H20O2	000103-09-3	22



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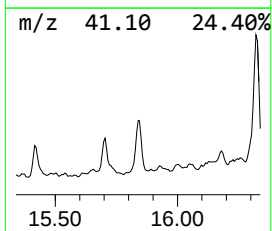
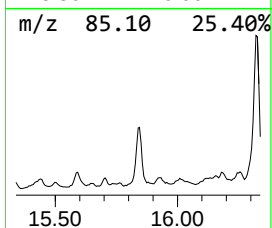
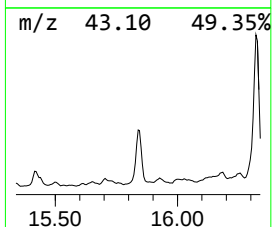
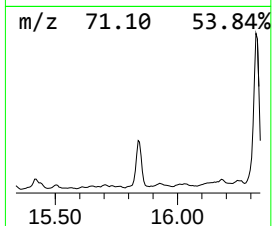
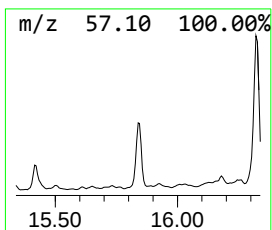
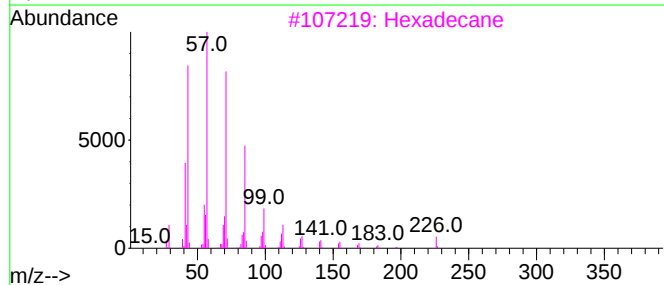
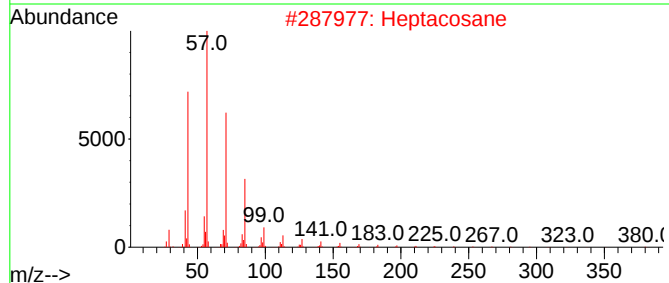
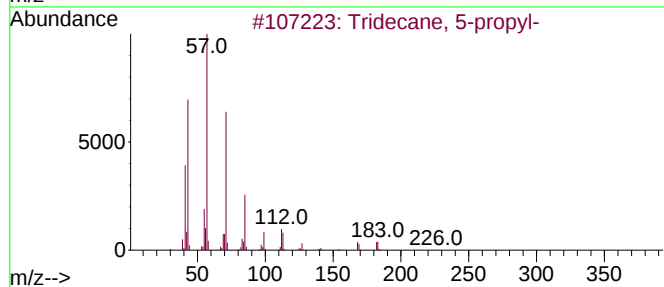
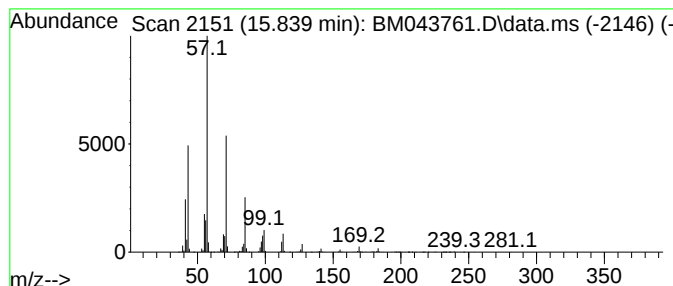
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.839	16.06 ng/ul	2220970	Acenaphthene-d10	14.598

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 5-propyl-	226	C16H34	055045-11-9	90
2			Heptacosane	380	C27H56	000593-49-7	90
3			Hexadecane	226	C16H34	000544-76-3	78
4			Hexadecane	226	C16H34	000544-76-3	72
5			Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
Data File : BM043761.D
Acq On : 05 Jan 2024 15:37
Operator : MA/JU
Sample : 05919-15ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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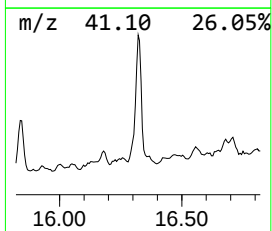
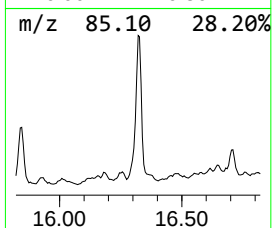
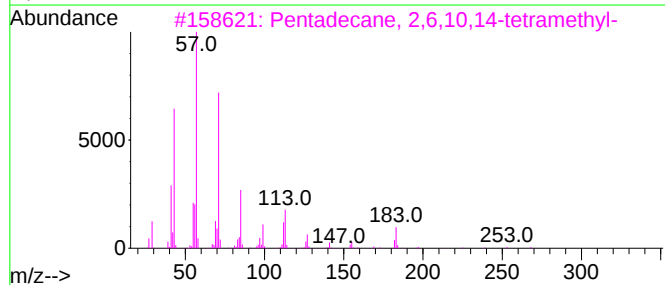
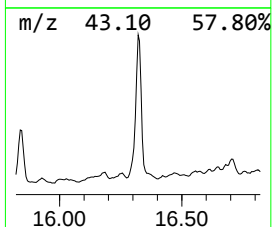
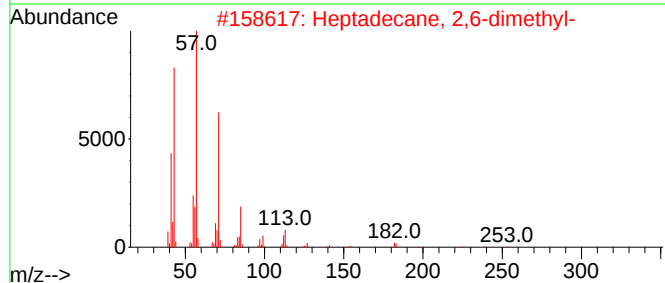
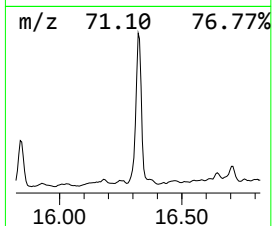
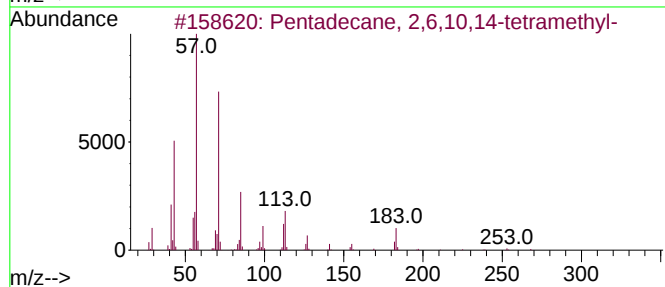
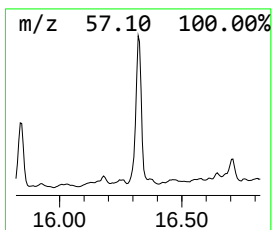
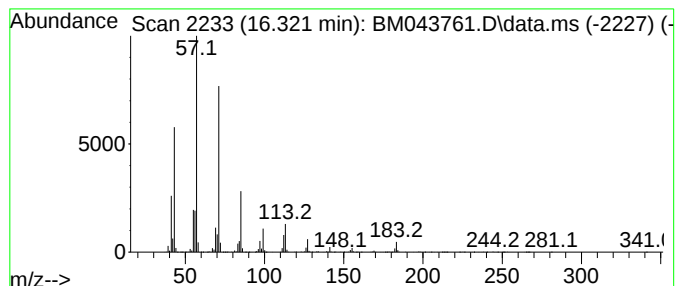
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.321	32.27 ng/ul	6436630	Phenanthrene-d10	17.356

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	91
2			Heptadecane, 2,6-dimethyl-	268	C19H40	054105-67-8	90
3			Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	90
4			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	86
5			Tridecane, 5-propyl-	226	C16H34	055045-11-9	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
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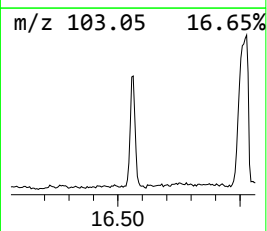
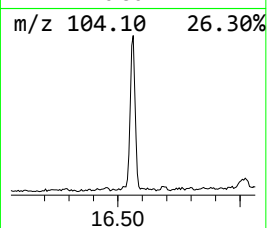
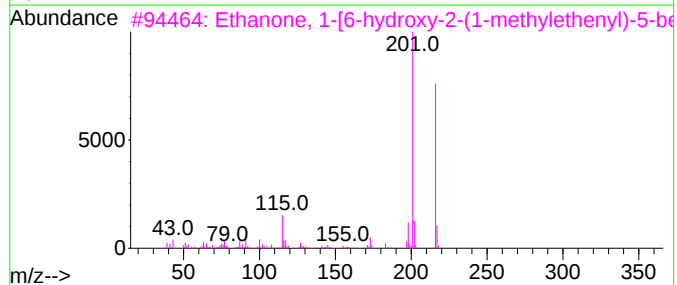
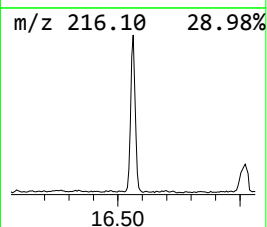
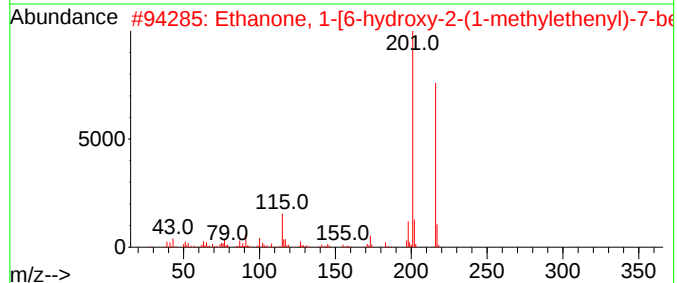
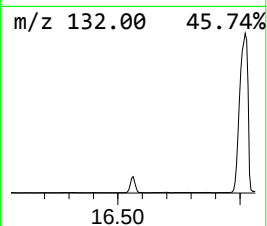
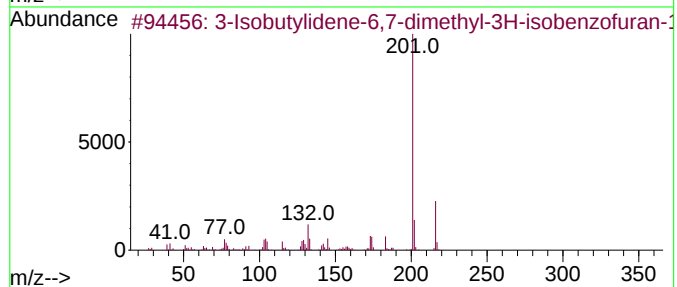
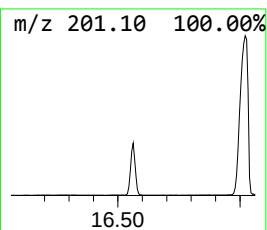
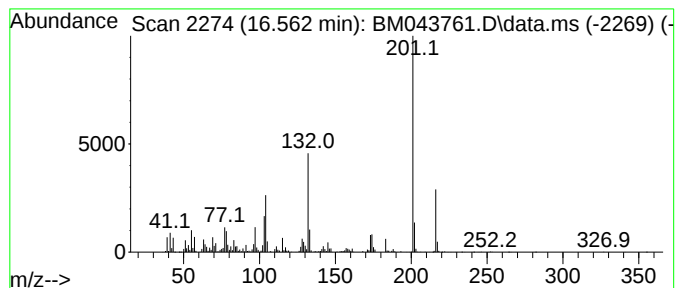
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 3-Isobutylidene-6,7-dimethy... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.562	6.86 ng/ul	1367480	Phenanthrene-d10	17.356

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Isobutylidene-6,7-dimethyl-3H-...	216	C14H16O2	1000186-84-7	52
2			Ethanone, 1-[6-hydroxy-2-(1-meth...	216	C13H12O3	055682-75-2	43
3			Ethanone, 1-[6-hydroxy-2-(1-meth...	216	C14H16O2	1000414-89-5	43
4			6-tert-Butyl-4-methylcoumarin	216	C14H16O2	017874-32-7	43
5			Leptaflorine	216	C13H16N2O	000486-93-1	43



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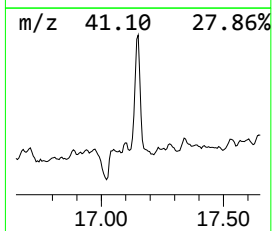
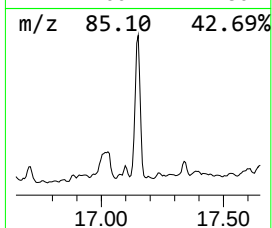
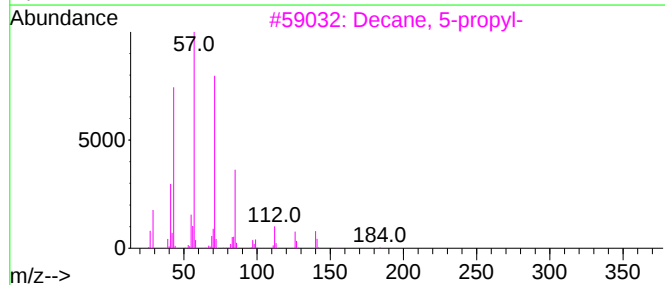
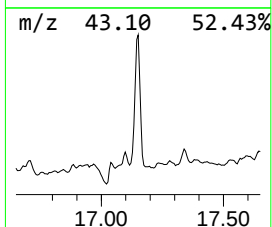
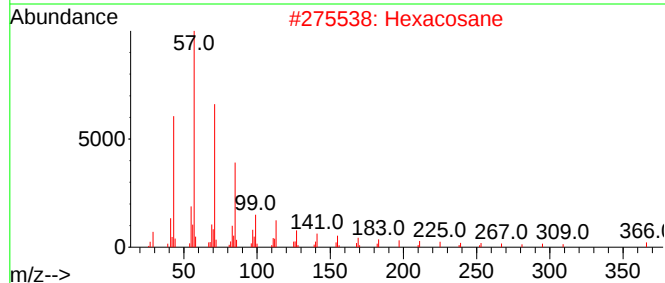
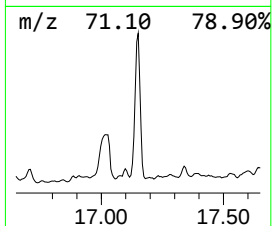
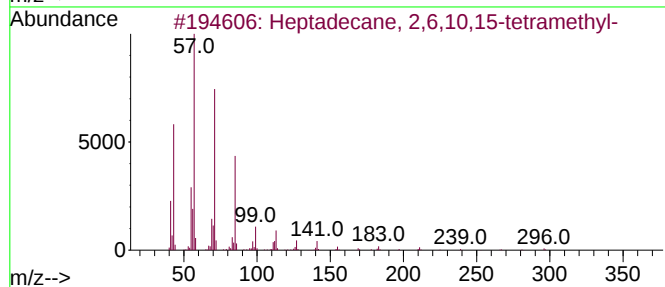
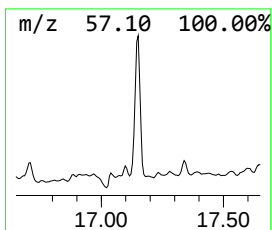
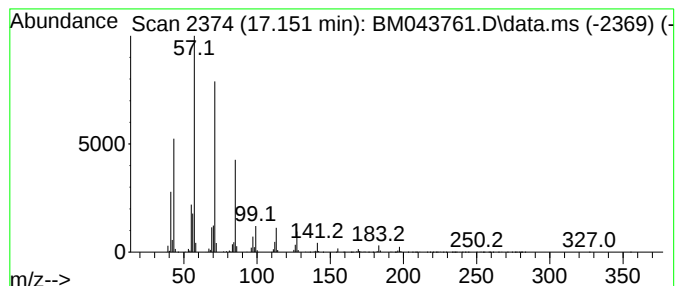
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.151	30.42 ng/ul	6066320	Phenanthrene-d10	17.356	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
2	Hexacosane	366	C26H54	000630-01-3	90
3	Decane, 5-propyl-	184	C13H28	017312-62-8	87
4	Decane, 3,7-dimethyl-	170	C12H26	017312-54-8	83
5	Dodecane, 4-methyl-	184	C13H28	006117-97-1	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
Data File : BM043761.D
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Instrument :
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ClientSampleId :
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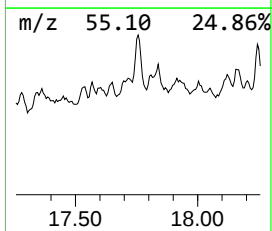
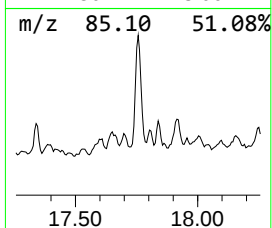
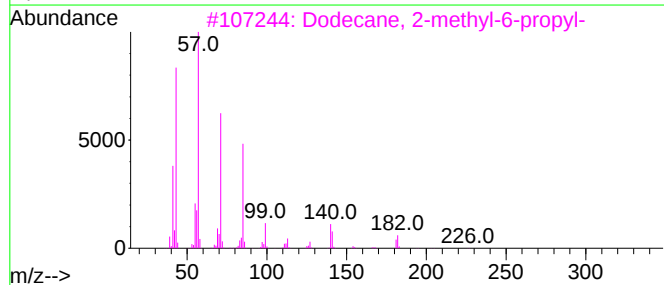
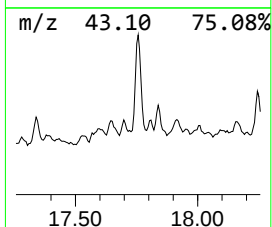
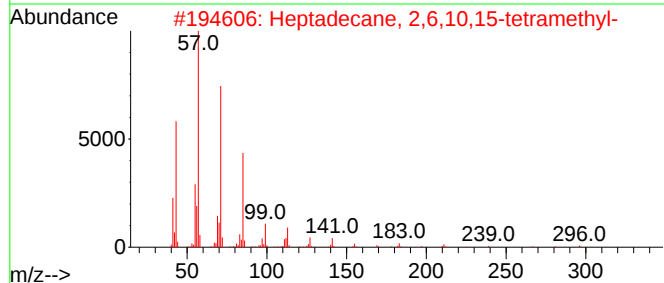
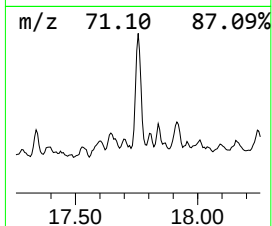
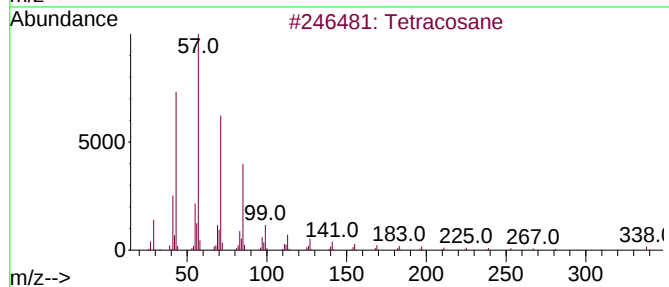
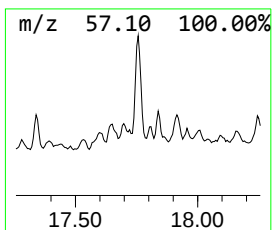
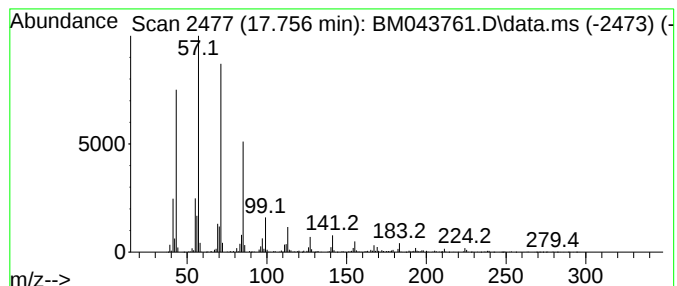
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.756	11.41 ng/ul	2276410	Phenanthrene-d10	17.356

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosane	338	C24H50	000646-31-1	91
2		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
3		Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	90
4		Heptadecane	240	C17H36	000629-78-7	90
5		Octadecane	254	C18H38	000593-45-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
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Sample : 05919-15ME
Misc :
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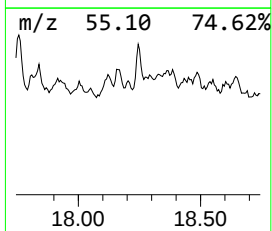
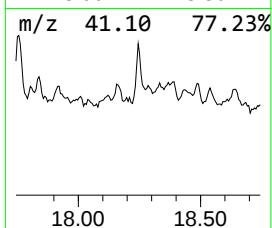
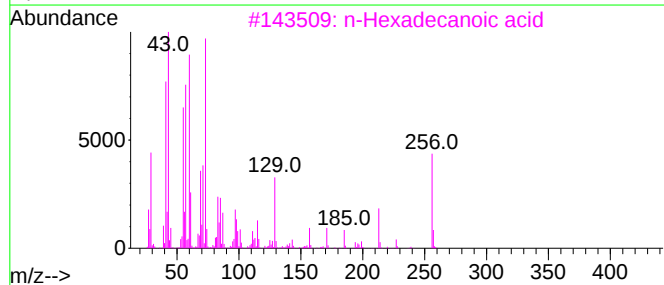
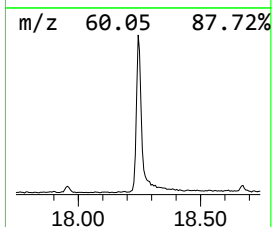
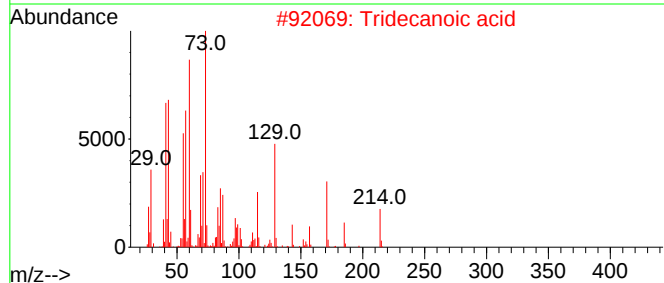
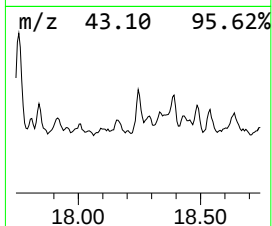
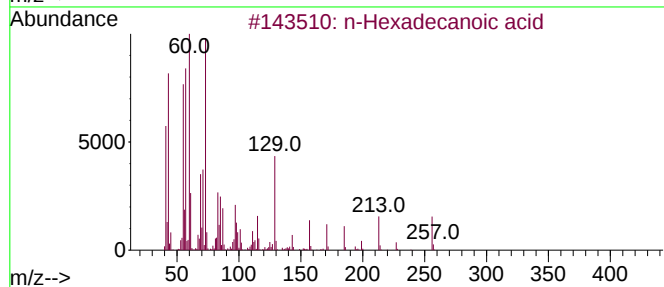
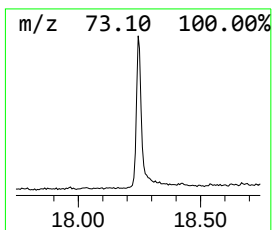
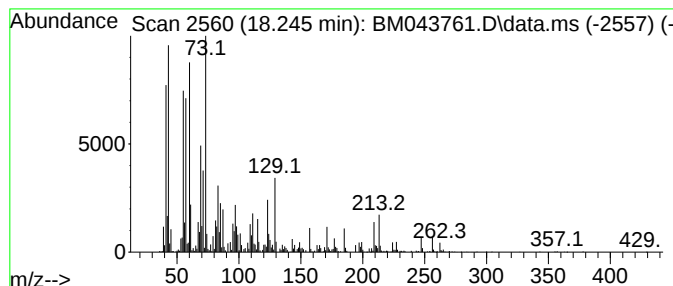
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.245	7.17 ng/ul	1430340	Phenanthrene-d10	17.356	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tridecanoic acid	214	C13H26O2	000638-53-9	96
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
5	Pentadecanoic acid	242	C15H30O2	001002-84-2	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
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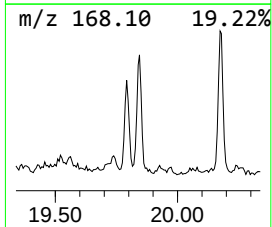
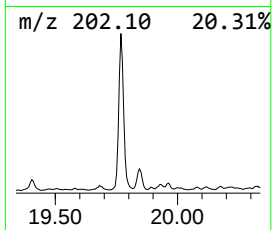
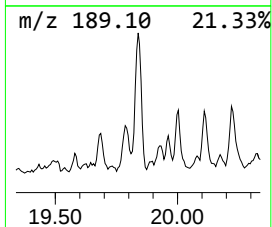
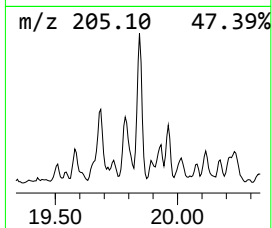
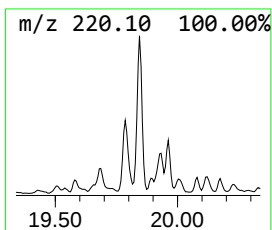
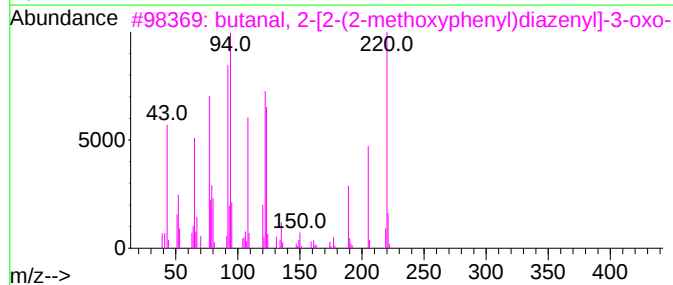
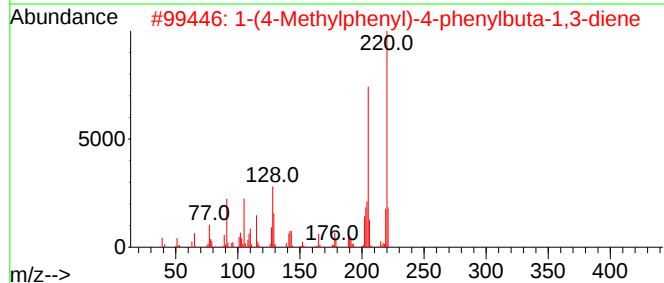
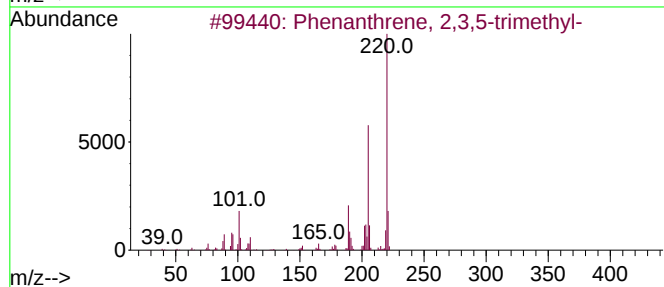
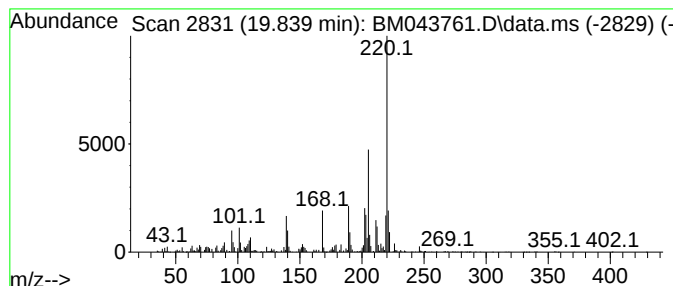
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Phenanthrene, 2,3,5-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.839	7.58 ng/ul	1416620	Chrysene-d12	21.533	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,3,5-trimethyl-	220	C17H16	003674-73-5	60
2	1-(4-Methylphenyl)-4-phenylbuta...	220	C17H16	037985-11-8	58
3	butanal, 2-[2-(2-methoxyphenyl)d...	220	C11H12N2O3	1000396-24-5	53
4	Phenanthrene, 2,3,5-trimethyl-	220	C17H16	003674-73-5	53
5	7-Nitro-6-methoxy-2,3-dimethylin...	220	C11H12N2O3	068289-71-4	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
Data File : BM043761.D
Acq On : 05 Jan 2024 15:37
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Misc :
ALS Vial : 6 Sample Multiplier: 1

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ClientSampleId :
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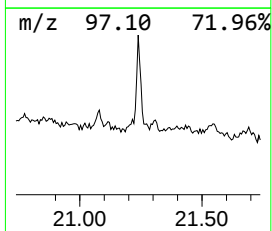
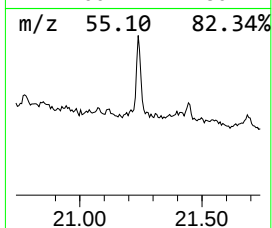
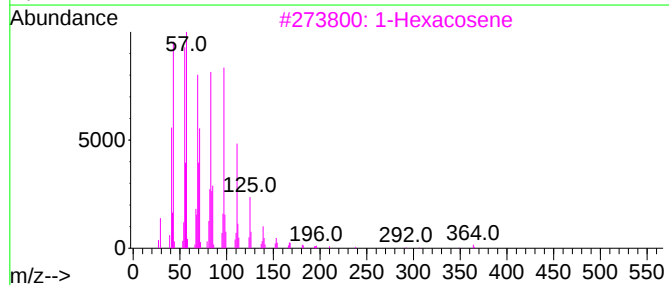
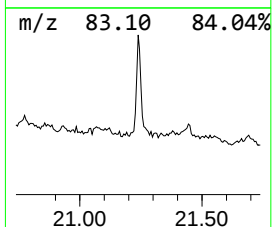
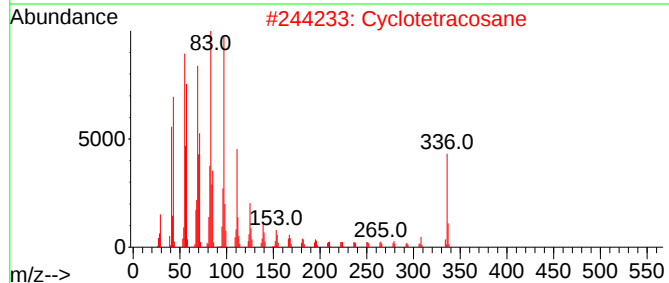
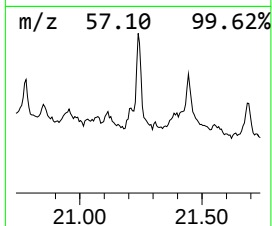
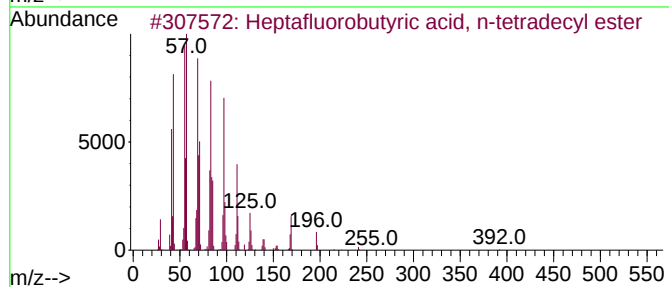
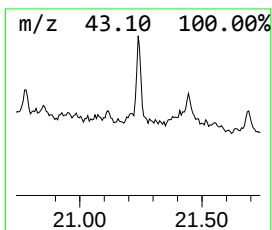
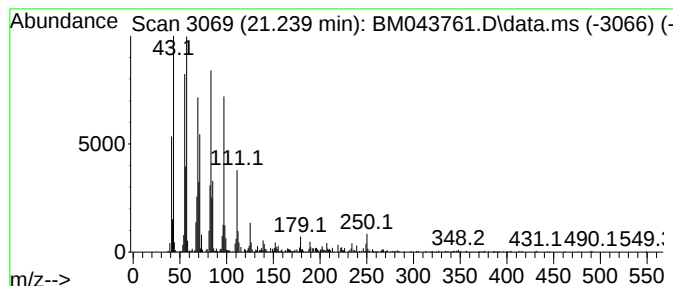
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Heptafluorobutyric acid, n-... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.238	4.62 ng/ul	863165	Chrysene-d12	21.533

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptafluorobutyric acid, n-tetra...	410	C18H29F7O2	007365-36-8	91
2			Cyclotetracosane	336	C24H48	000297-03-0	91
3			1-Hexacosene	364	C26H52	018835-33-1	91
4			Pentacos-1-ene	350	C25H50	016980-85-1	91
5			n-Nonadecanol-1	284	C19H40O	001454-84-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
Data File : BM043761.D
Acq On : 05 Jan 2024 15:37
Operator : MA/JU
Sample : 05919-15ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCF47ME

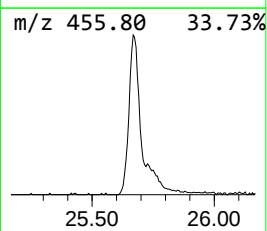
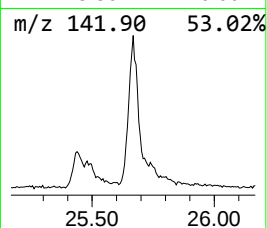
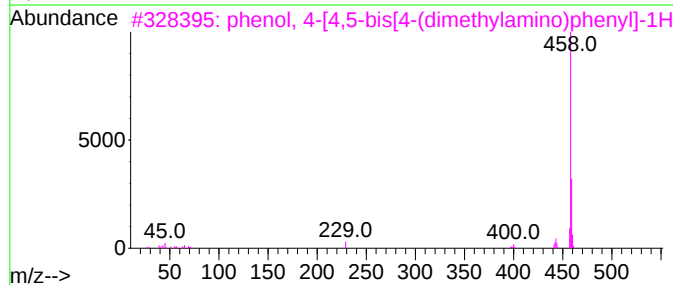
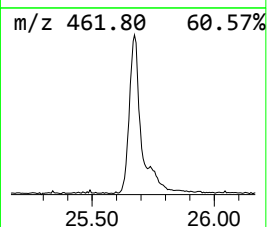
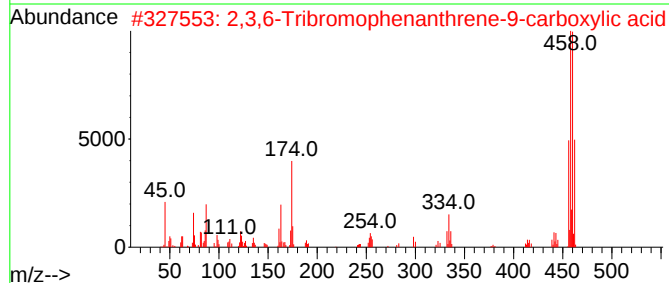
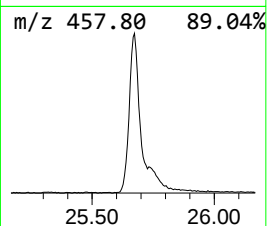
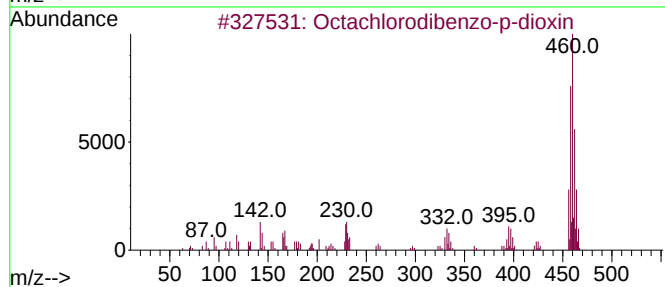
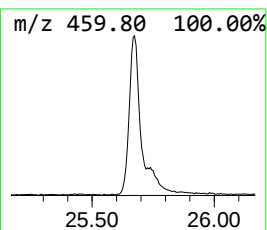
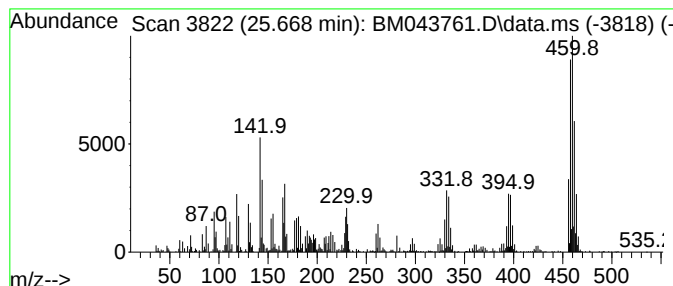
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Octachlorodibenzo-p-dioxin Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.668	9.70 ng/ul	1778290	Perylene-d12	23.950

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octachlorodibenzo-p-dioxin	456	C12Cl8O2	003268-87-9	90
2			2,3,6-Tribromophenanthrene-9-car...	456	C15H7Br3O2	056988-60-4	10
3			phenol, 4-[4,5-bis[4-(dimethylam...	458	C27H30N4O3	1000399-43-0	9
4			1,1':4',1'':4'',1''':4''',1''''...	458	C36H26	004499-83-6	9
5			1,1':4',1'':4'',1''':4''',1''''...	458	C36H26	004499-83-6	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
Data File : BM043761.D
Acq On : 05 Jan 2024 15:37
Operator : MA/JU
Sample : 05919-15ME
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCF47ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.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
(DEL) Alkane: S...	14.068	3.8	ng/ul	525003	3	14.598	2765420	20.0
2-Ethylhexyl me...	14.480	26.1	ng/ul	3610530	3	14.598	2765420	20.0
(DEL) Alkane: S...	15.104	2.8	ng/ul	391019	3	14.598	2765420	20.0
2-Ethylhexyl 2-...	15.421	9.9	ng/ul	1371370	3	14.598	2765420	20.0
(DEL) Alkane: S...	15.839	16.1	ng/ul	2220970	3	14.598	2765420	20.0
(DEL) Alkane: S...	16.321	32.3	ng/ul	6436630	4	17.356	3988730	20.0
3-Isobutylidene...	16.562	6.9	ng/ul	1367480	4	17.356	3988730	20.0
(DEL) Alkane: S...	17.151	30.4	ng/ul	6066320	4	17.356	3988730	20.0
(DEL) Alkane: S...	17.756	11.4	ng/ul	2276410	4	17.356	3988730	20.0
n-Hexadecanoic ...	18.245	7.2	ng/ul	1430340	4	17.356	3988730	20.0
Phenanthrene, 2...	19.839	7.6	ng/ul	1416620	5	21.533	3738240	20.0
Heptafluorobuty...	21.238	4.6	ng/ul	863165	5	21.533	3738240	20.0
Octachlorodiben...	25.668	9.7	ng/ul	1778290	6	23.950	3667970	20.0