

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017563.D
 Acq On : 05 Oct 2023 13:40
 Operator : MA/JU
 Sample : 04679-04
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PMW-4(20230927)

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-BP100423.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP017563.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.287	31	34	38	rBV2	21179	25262	0.48%	0.057%
2	3.599	84	87	91	rBV5	5685	8282	0.16%	0.019%
3	3.728	105	109	120	rBV	191138	289812	5.54%	0.652%
4	3.916	138	141	144	rVB4	6518	7275	0.14%	0.016%
5	4.028	157	160	164	rVB	10134	12741	0.24%	0.029%
6	5.193	353	358	367	rVB	37817	57012	1.09%	0.128%
7	6.381	556	560	564	rBV7	2963	5721	0.11%	0.013%
8	6.905	648	649	654	rVB5	5461	5556	0.11%	0.013%
9	7.099	675	682	697	rBV	150909	267812	5.12%	0.603%
10	7.281	706	713	723	rBV	522723	798772	15.27%	1.798%
11	7.457	736	743	754	rBV	537957	840366	16.06%	1.892%
12	7.552	754	759	762	rVV3	4635	6963	0.13%	0.016%
13	7.810	798	803	810	rBV	119268	192685	3.68%	0.434%
14	7.940	818	825	829	rBV	412604	701279	13.40%	1.579%
15	7.975	829	831	839	rVB	134202	165850	3.17%	0.373%
16	8.169	858	864	868	rBV5	6662	11605	0.22%	0.026%
17	8.287	880	884	890	rVB	14827	23742	0.45%	0.053%
18	8.657	938	947	958	rBV	485335	805112	15.39%	1.812%
19	9.146	1023	1030	1042	rBV	619126	1016695	19.43%	2.289%
20	9.257	1045	1049	1055	rVB9	3672	6957	0.13%	0.016%
21	9.340	1060	1063	1070	rVB4	9272	14274	0.27%	0.032%
22	9.863	1145	1152	1165	rBV	522101	871243	16.65%	1.961%
23	10.057	1177	1185	1187	rBV9	4939	9573	0.18%	0.022%
24	10.393	1233	1242	1256	rBV	745005	1337625	25.57%	3.011%
25	10.622	1274	1281	1289	rVB	232167	384104	7.34%	0.865%
26	10.775	1299	1307	1315	rBV	574661	939046	17.95%	2.114%
27	10.945	1329	1336	1348	rBV	607357	1050155	20.07%	2.364%
28	11.145	1365	1370	1376	rVB10	4740	9186	0.18%	0.021%
29	11.469	1419	1425	1428	rBV3	8948	12485	0.24%	0.028%
30	11.551	1432	1439	1443	rVB9	4393	7458	0.14%	0.017%
31	11.804	1474	1482	1489	rBV5	24809	62245	1.19%	0.140%
32	12.316	1563	1569	1571	rBV7	3347	5539	0.11%	0.012%
33	12.369	1572	1578	1585	rVB2	13989	25023	0.48%	0.056%
34	12.669	1626	1629	1634	rVB7	4638	8127	0.16%	0.018%
35	13.469	1760	1765	1775	rVB2	19001	32587	0.62%	0.073%
36	13.804	1817	1822	1823	rBV5	3558	5899	0.11%	0.013%

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37	14.045	1857	1863	1875	rBV	1559436	2099721	40.14%	4.727%
38	14.245	1892	1897	1903	rVB8	15239	24811	0.47%	0.056%
39	14.328	1903	1911	1922	rBV	1652186	2405329	45.98%	5.415%
40	14.628	1956	1962	1970	rBV	885358	1269541	24.27%	2.858%
41	14.886	2000	2006	2019	rBV	84939	175124	3.35%	0.394%
42	15.016	2025	2028	2031	rBV4	23434	29572	0.57%	0.067%
43	15.051	2031	2034	2038	rVB6	10305	13373	0.26%	0.030%
44	15.375	2085	2089	2094	rVB	21105	29232	0.56%	0.066%
45	15.633	2126	2133	2143	rBV	2308468	3135658	59.94%	7.059%
46	15.775	2152	2157	2167	rBV	725341	985061	18.83%	2.217%
47	15.869	2170	2173	2178	rVB6	12324	21281	0.41%	0.048%
48	16.151	2217	2221	2228	rVB	23211	31068	0.59%	0.070%
49	16.769	2323	2326	2331	rVB7	11263	13805	0.26%	0.031%
50	17.422	2431	2437	2447	rBV	1156508	1639592	31.34%	3.691%
51	17.522	2448	2454	2466	rVV	2598710	3488934	66.69%	7.854%
52	18.269	2576	2581	2591	rBV	267663	402418	7.69%	0.906%
53	18.692	2650	2653	2660	rVB	33946	51205	0.98%	0.115%
54	19.345	2761	2764	2767	rBV	29612	33053	0.63%	0.074%
55	19.416	2771	2776	2782	rVB	338945	418098	7.99%	0.941%
56	19.510	2789	2792	2801	rBV	126078	207326	3.96%	0.467%
57	19.651	2811	2816	2821	rBV	1105542	1180906	22.57%	2.658%
58	19.710	2824	2826	2830	rVB	39837	40886	0.78%	0.092%
59	19.774	2831	2837	2843	rBV	3615001	4672830	89.32%	10.519%
60	20.474	2953	2956	2960	rVB	187470	192908	3.69%	0.434%
61	20.680	2987	2991	2996	rBV	847590	897662	17.16%	2.021%
62	20.733	2997	3000	3005	rVV	98318	120661	2.31%	0.272%
63	20.780	3005	3008	3011	rVB3	62484	75610	1.45%	0.170%
64	21.557	3135	3140	3147	rBV	1528862	1953859	37.35%	4.398%
65	21.651	3153	3156	3160	rVB	136681	147824	2.83%	0.333%
66	22.133	3235	3238	3244	rVB2	136719	163480	3.12%	0.368%
67	22.663	3325	3328	3333	rVB	188910	250754	4.79%	0.564%
68	23.262	3426	3430	3436	rVB	237619	373827	7.15%	0.842%
69	23.968	3541	3550	3561	rVB2	2290057	5231565	100.00%	11.777%
70	24.139	3572	3579	3589	rVB	965155	2082070	39.80%	4.687%
71	24.739	3676	3681	3691	rVB	251858	539202	10.31%	1.214%

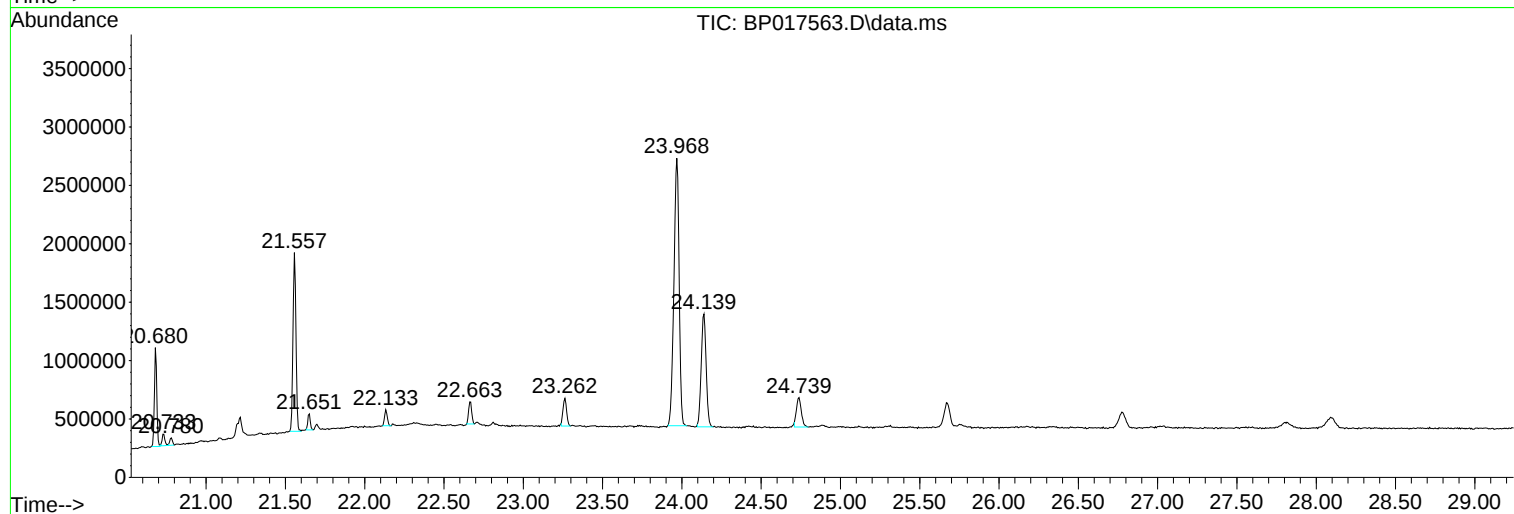
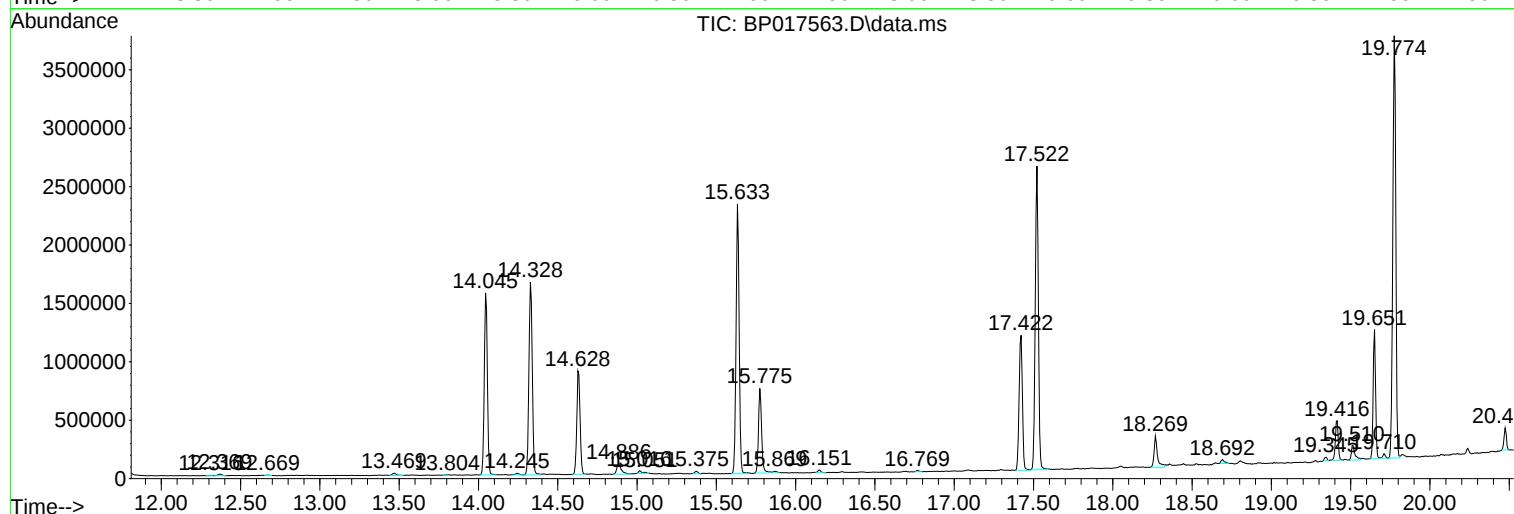
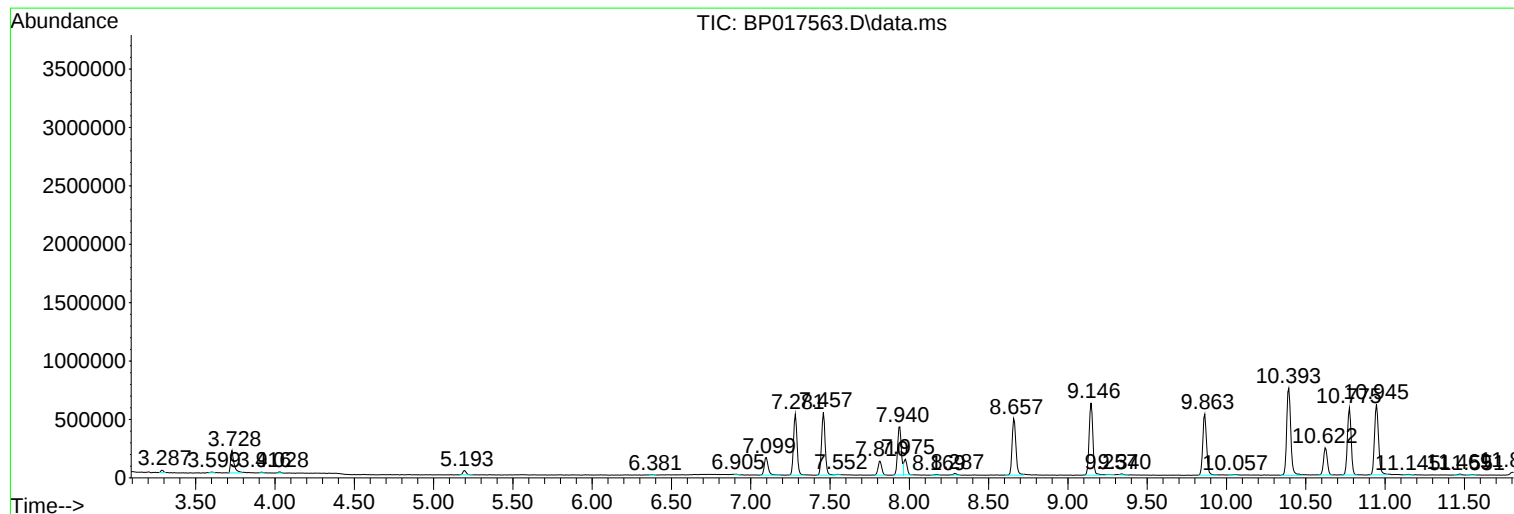
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ClientSampleId :
PMW-4(20230927)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.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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PMW-4(20230927)

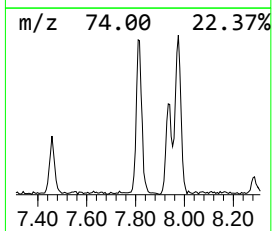
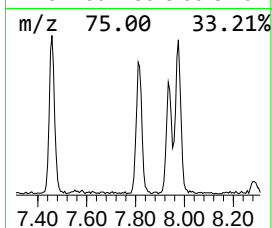
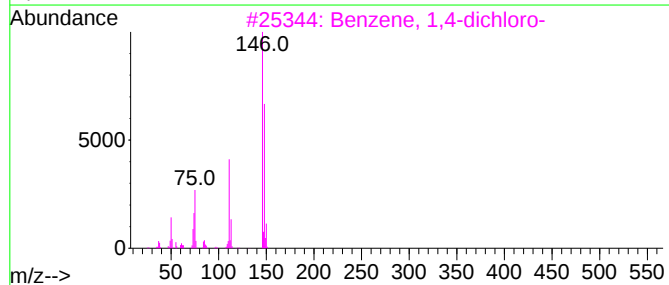
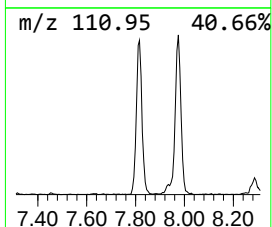
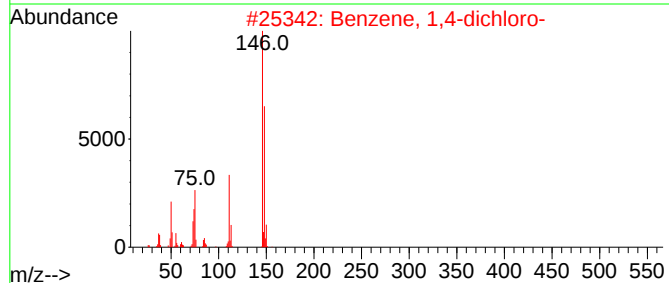
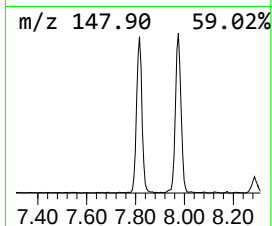
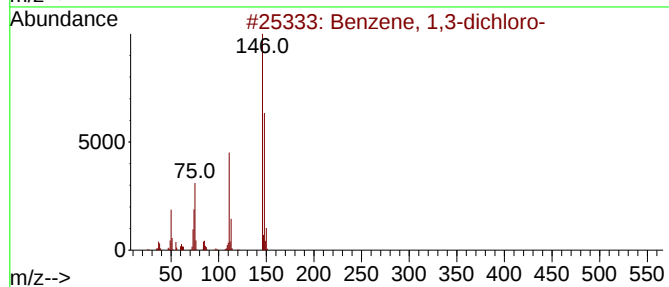
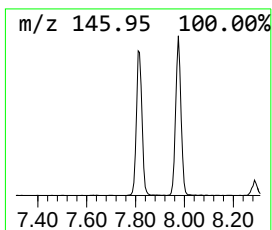
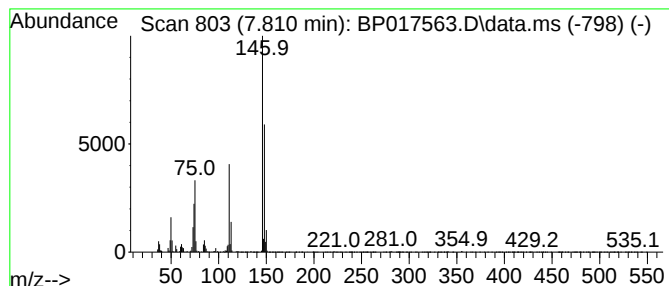
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, 1,3-dichloro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.810	5.50 ng/ul	192685	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	98
2			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	97
3			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	96
4			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96
5			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	96



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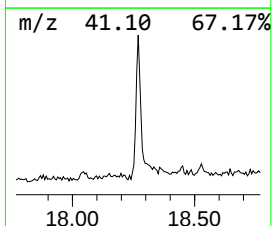
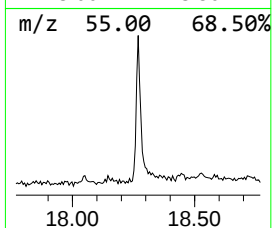
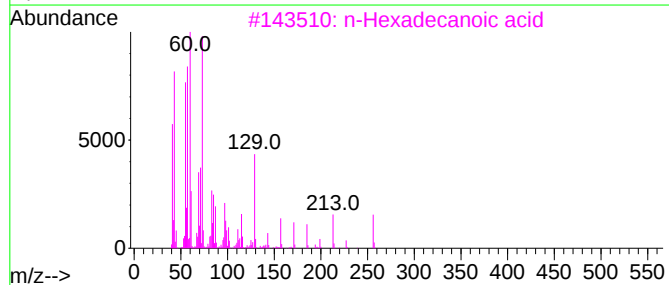
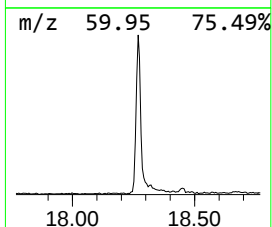
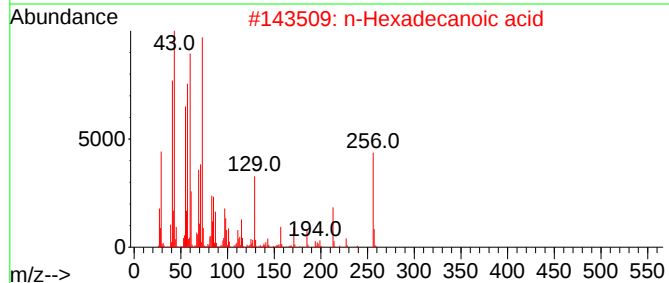
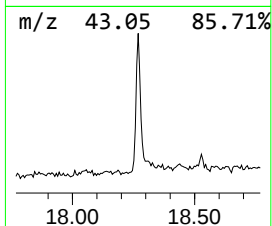
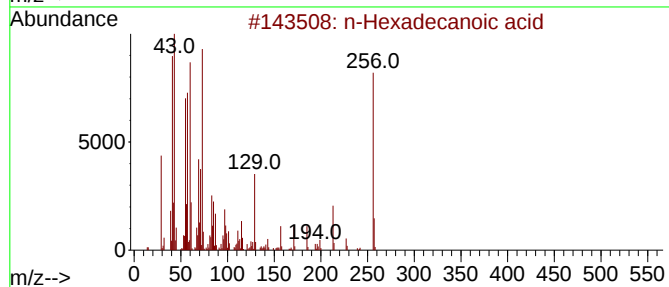
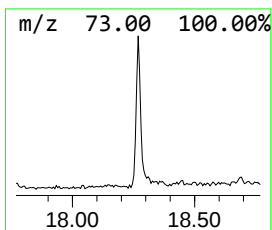
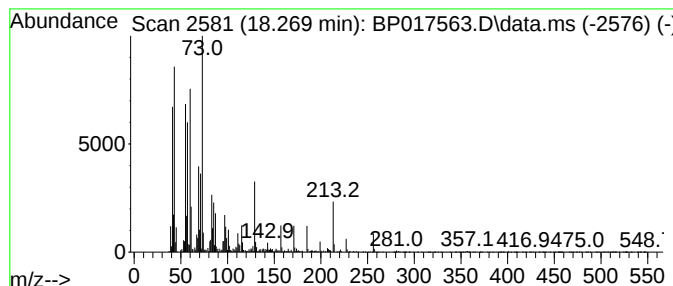
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.269	4.91 ng/ul	402418	Phenanthrene-d10	17.422

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	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	90
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	86



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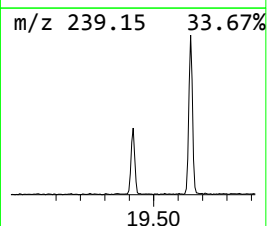
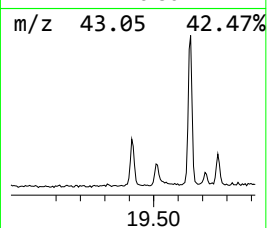
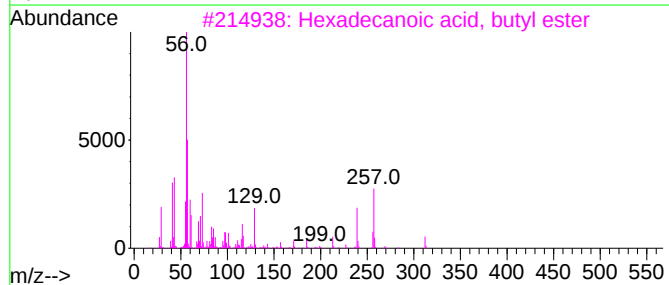
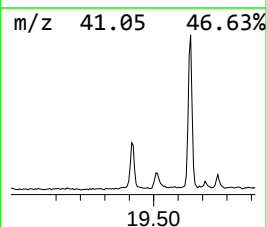
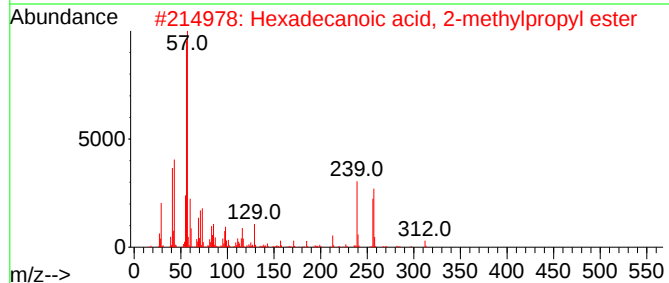
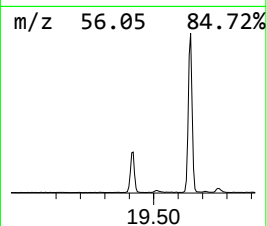
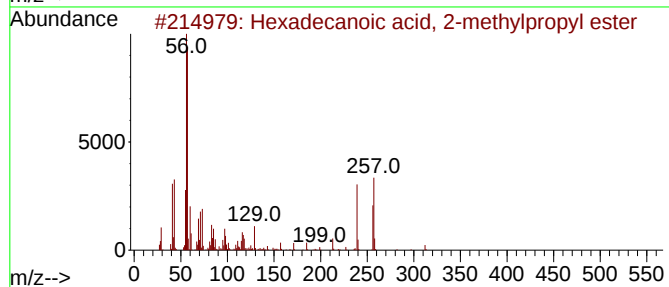
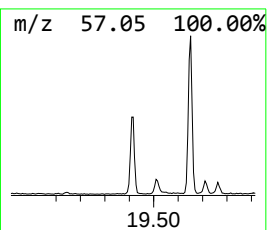
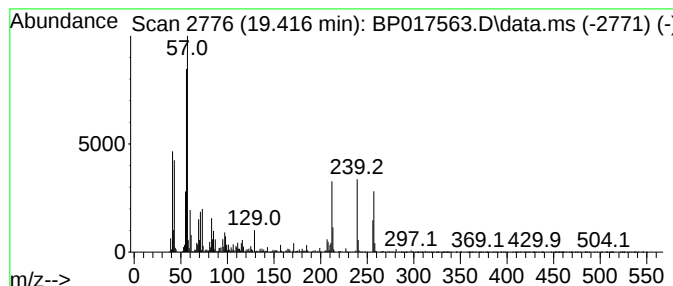
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Peak Number 4 Hexadecanoic acid, 2-methyl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.416	5.10 ng/ul	418098	Phenanthrene-d10	17.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanoic acid, 2-methylpropy...	312	C20H40O2	000110-34-9	74
2			Hexadecanoic acid, 2-methylpropy...	312	C20H40O2	000110-34-9	68
3			Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	53
4			Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	53
5			Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	27



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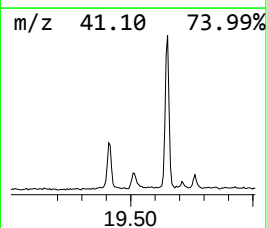
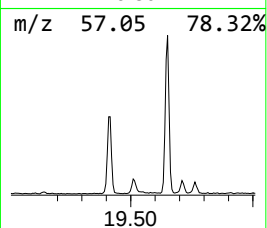
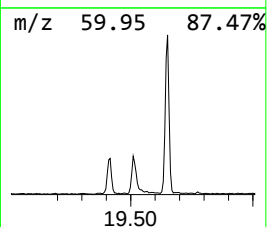
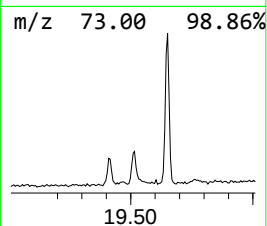
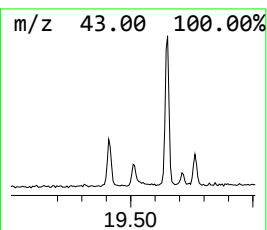
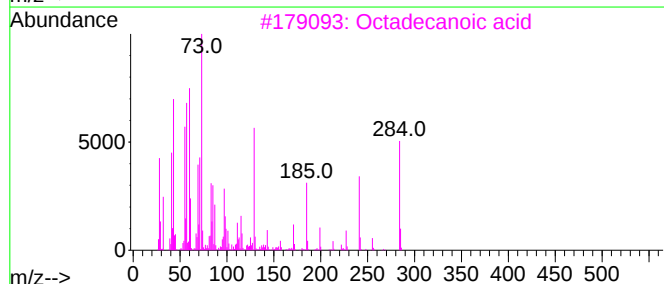
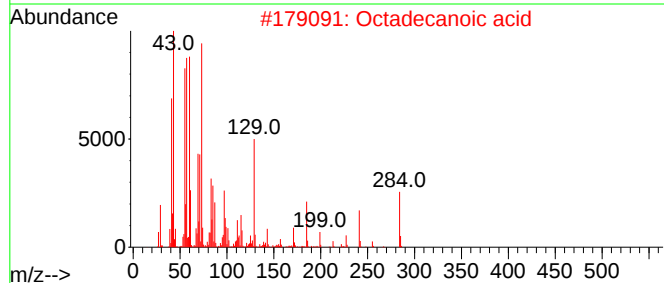
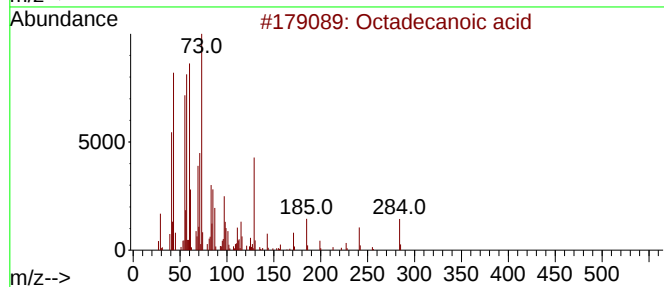
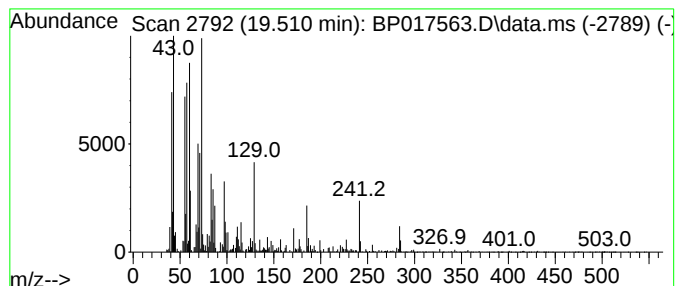
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.510	2.12 ng/ul	207326	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	98
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\BP100423\
Data File : BP017563.D
Acq On : 05 Oct 2023 13:40
Operator : MA/JU
Sample : 04679-04
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PMW-4(20230927)

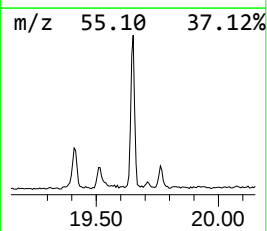
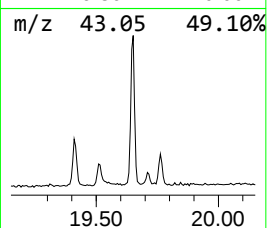
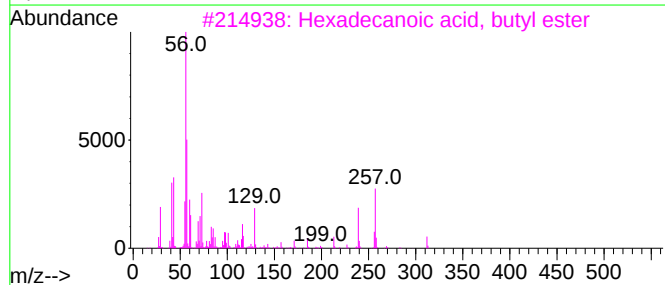
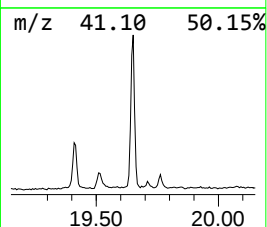
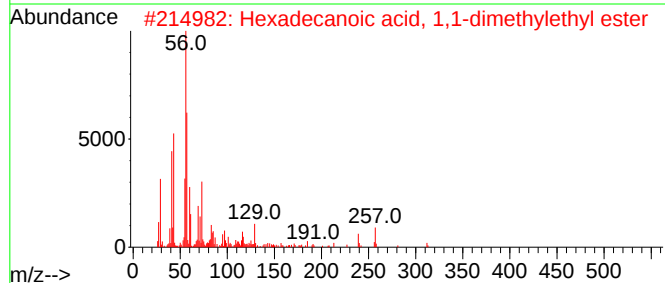
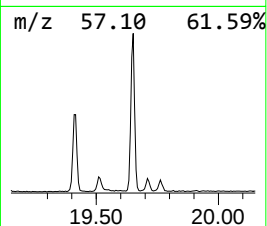
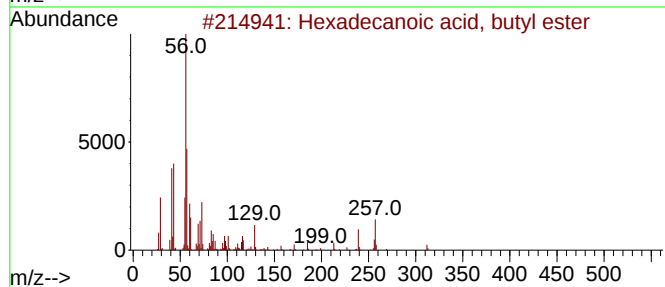
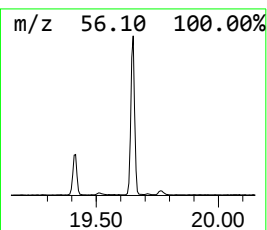
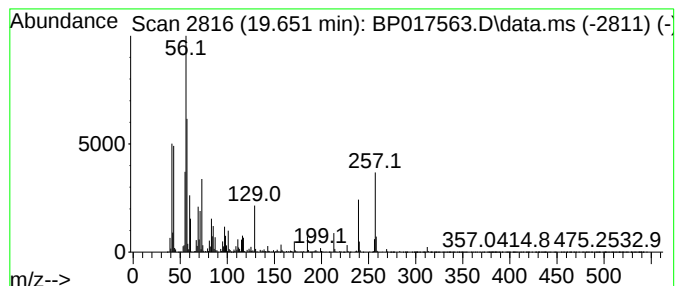
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Hexadecanoic acid, butyl ester Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.651	12.09 ng/ul	1180910	Chrysene-d12	21.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	97
2			Hexadecanoic acid, 1,1-dimethyle...	312	C20H40O2	031158-91-5	96
3			Hexadecanoic acid, butyl ester	312	C20H40O2	000111-06-8	91
4			Hexadecanoic acid, 2-methylpropy...	312	C20H40O2	000110-34-9	78
5			Hexadecanoic acid, 2-methylpropy...	312	C20H40O2	000110-34-9	45



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017563.D
Acq On : 05 Oct 2023 13:40
Operator : MA/JU
Sample : 04679-04
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PMW-4(20230927)

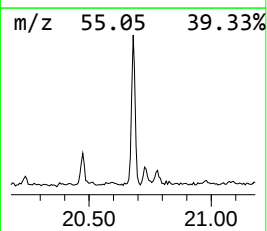
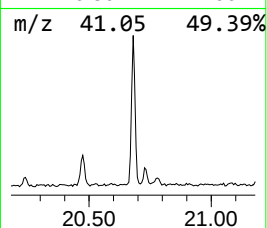
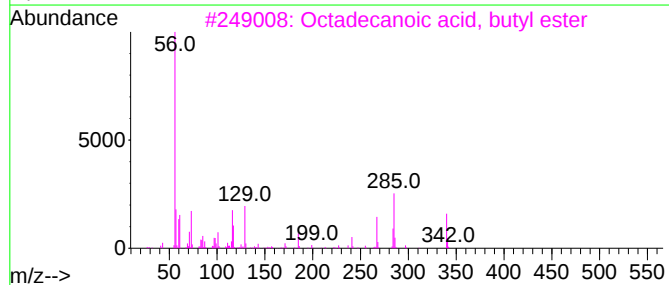
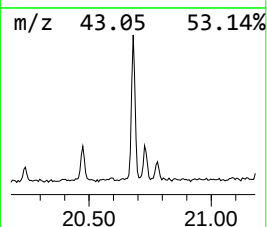
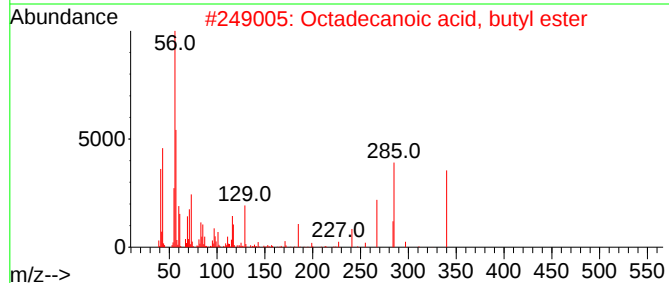
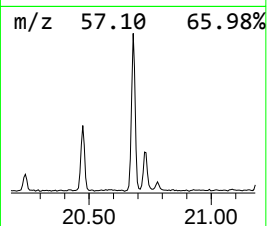
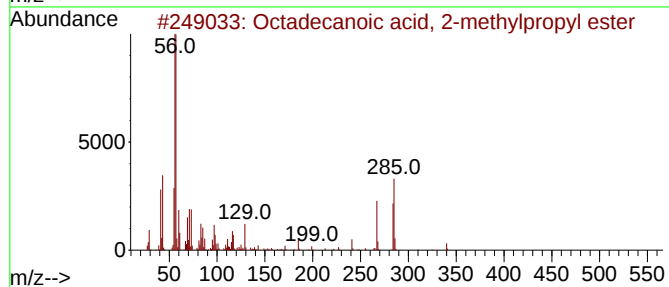
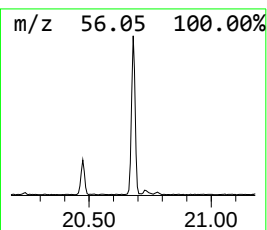
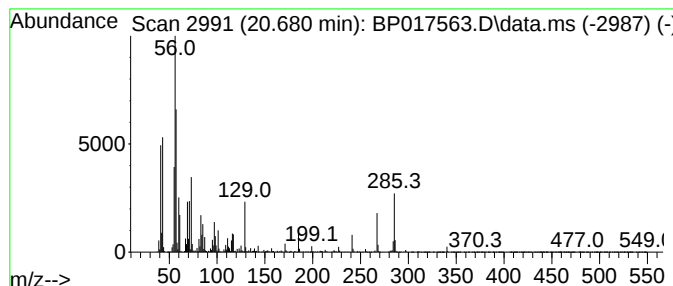
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Octadecanoic acid, 2-methyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.680	9.19 ng/ul	897662	Chrysene-d12	21.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid, 2-methylpropy...	340	C22H44O2	000646-13-9	89
2			Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	87
3			Octadecanoic acid, butyl ester	340	C22H44O2	000123-95-5	80
4			Octadecanoic acid, 2-methylpropy...	340	C22H44O2	000646-13-9	74
5			Butyl myristate	284	C18H36O2	000110-36-1	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017563.D
Acq On : 05 Oct 2023 13:40
Operator : MA/JU
Sample : 04679-04
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
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ClientSampleId :
PMW-4(20230927)

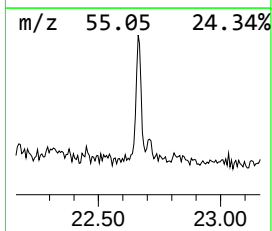
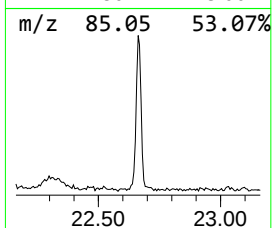
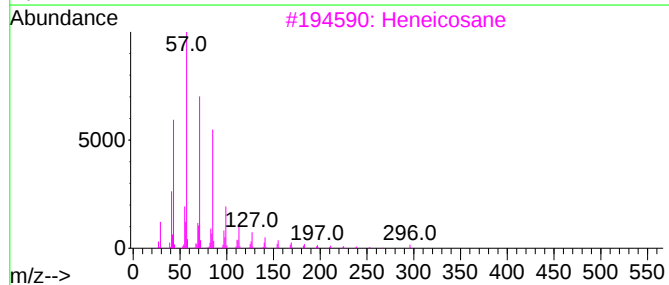
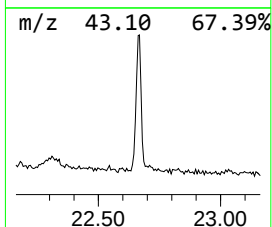
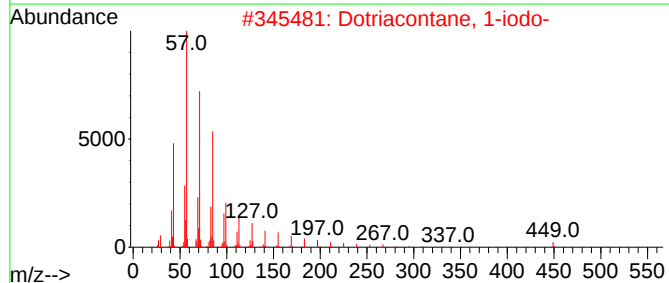
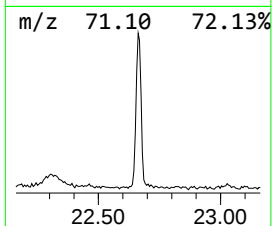
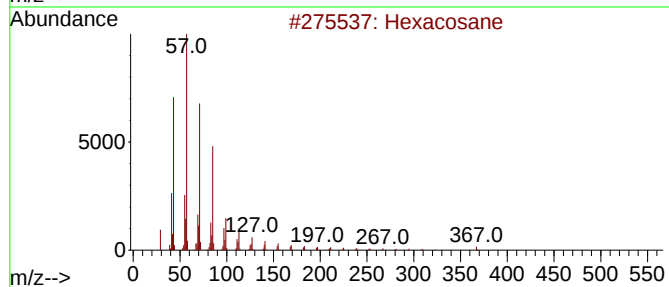
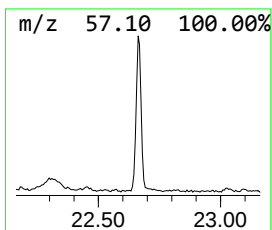
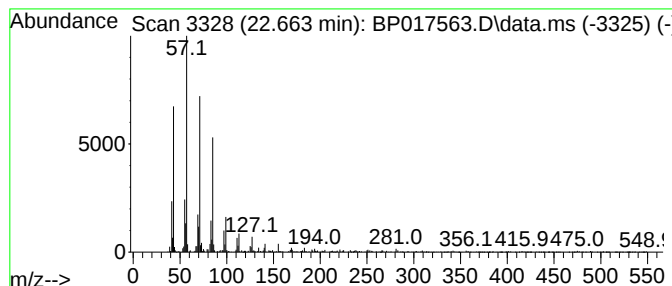
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.663	2.57 ng/ul	250754	Chrysene-d12	21.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	90
2			Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	87
3			Heneicosane	296	C21H44	000629-94-7	87
4			3-Methylheptacosane	394	C28H58	014167-66-9	86
5			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017563.D
Acq On : 05 Oct 2023 13:40
Operator : MA/JU
Sample : 04679-04
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PMW-4(20230927)

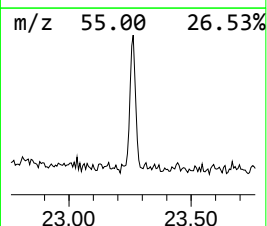
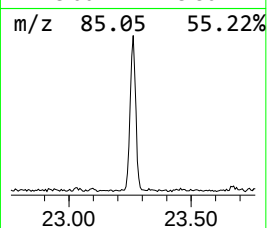
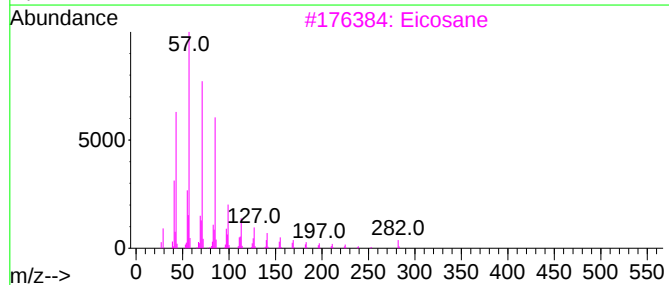
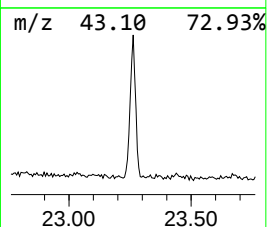
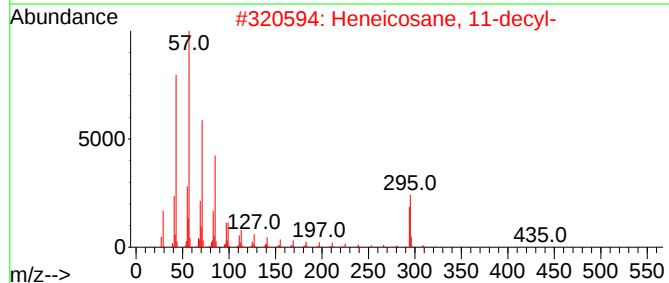
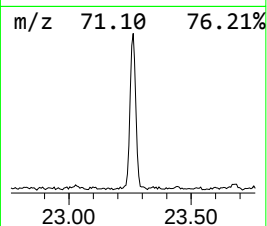
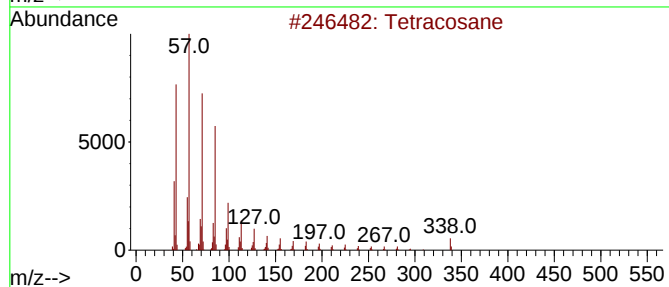
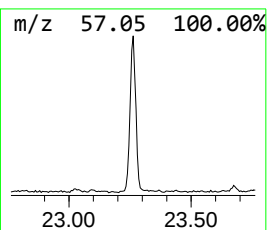
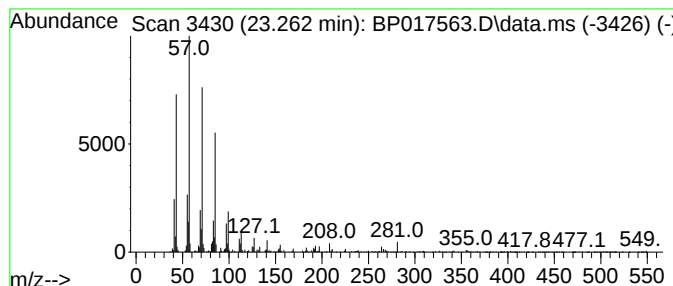
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.262	3.59 ng/ul	373827	Perylene-d12	24.139

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosane	338	C24H50	000646-31-1	96
2		Heneicosane, 11-decyl-	437	C31H64	055320-06-4	95
3		Eicosane	282	C20H42	000112-95-8	93
4		Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
5		Octacosane, 2-methyl-	408	C29H60	001560-98-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017563.D
Acq On : 05 Oct 2023 13:40
Operator : MA/JU
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Misc :
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ClientSampleId :
PMW-4(20230927)

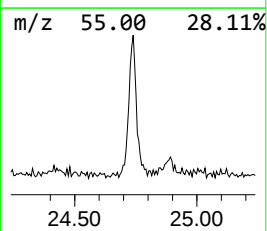
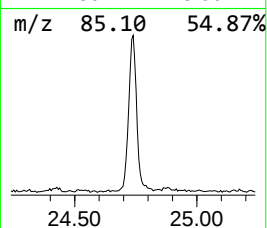
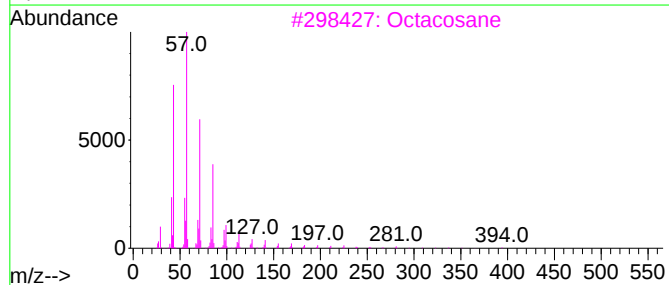
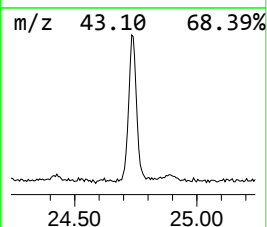
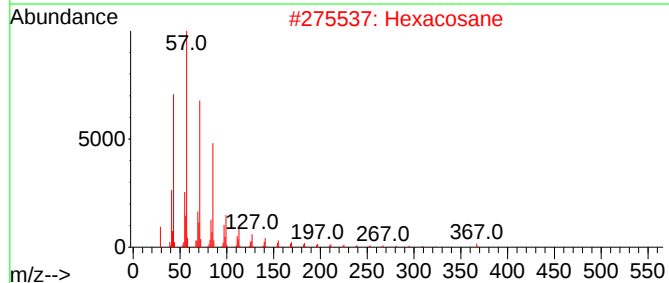
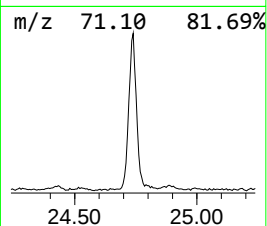
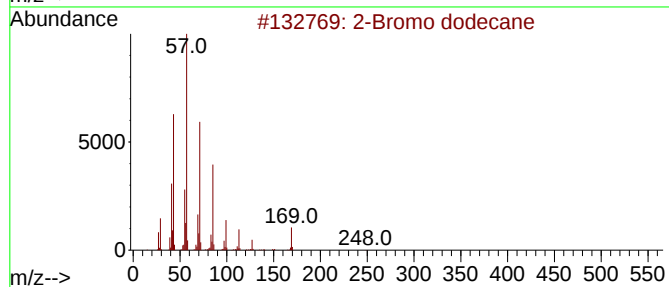
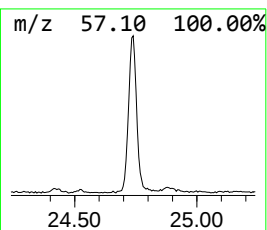
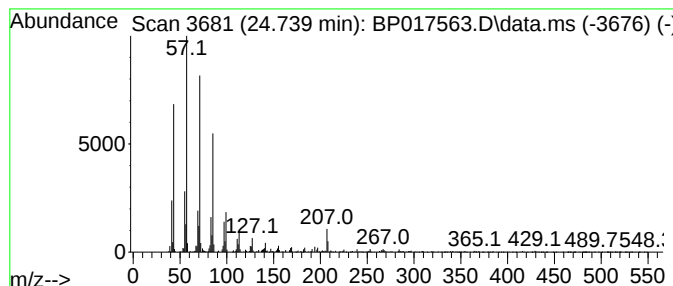
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 2-Bromo dodecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.739	5.18 ng/ul	539202	Perylene-d12	24.139

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Bromo dodecane	248	C12H25Br	013187-99-0	91
2		Hexacosane	366	C26H54	000630-01-3	90
3		Octacosane	394	C28H58	000630-02-4	87
4		Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	87
5		Octadecane	254	C18H38	000593-45-3	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017563.D
Acq On : 05 Oct 2023 13:40
Operator : MA/JU
Sample : 04679-04
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PMW-4(20230927)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.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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n-Hexadecanoic ...	18.269	4.9	ng/ul	402418	4	17.422	1639590	20.0
Hexadecanoic ac...	19.416	5.1	ng/ul	418098	4	17.422	1639590	20.0
Octadecanoic acid	19.510	2.1	ng/ul	207326	5	21.557	1953860	20.0
Hexadecanoic ac...	19.651	12.1	ng/ul	1180910	5	21.557	1953860	20.0
Octadecanoic ac...	20.680	9.2	ng/ul	897662	5	21.557	1953860	20.0
(DEL) Alkane: S...	22.663	2.6	ng/ul	250754	5	21.557	1953860	20.0
(DEL) Alkane: S...	23.262	3.6	ng/ul	373827	6	24.139	2082070	20.0
2-Bromo dodecane	24.739	5.2	ng/ul	539202	6	24.139	2082070	20.0