

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
 Data File : BP013328.D
 Acq On : 25 Dec 2022 08:09
 Operator : CG/JU
 Sample : N6018-19
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBYK4

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-BP120722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP013328.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.328	20	24	30	rBV	151923	188744	1.36%	0.133%
2	3.758	93	97	111	rBV	1758791	2625036	18.86%	1.855%
3	4.081	149	152	156	rBV	40059	53856	0.39%	0.038%
4	5.028	309	313	320	rVB2	38882	61468	0.44%	0.043%
5	5.258	350	352	360	rVB7	9720	16896	0.12%	0.012%
6	6.605	578	581	588	rVB2	25295	42457	0.31%	0.030%
7	6.916	628	634	641	rBV10	10223	23502	0.17%	0.017%
8	7.099	657	665	676	rBV	2142202	3457628	24.84%	2.444%
9	7.263	686	693	704	rBV	1793188	2702567	19.42%	1.910%
10	7.369	708	711	719	rVB8	14662	27855	0.20%	0.020%
11	7.469	721	728	738	rVB	2206718	3465993	24.90%	2.450%
12	7.940	797	808	815	rBV	2100981	3354558	24.10%	2.371%
13	8.652	919	929	941	rBV	2400532	3931507	28.25%	2.779%
14	9.040	990	995	999	rBV2	31716	52277	0.38%	0.037%
15	9.105	999	1006	1022	rVV	1982005	3197846	22.98%	2.260%
16	9.263	1029	1033	1039	rVB9	15545	26265	0.19%	0.019%
17	9.505	1066	1074	1079	rBV2	42787	71881	0.52%	0.051%
18	9.693	1101	1106	1111	rVB	20601	31765	0.23%	0.022%
19	9.828	1122	1129	1139	rBV	1624303	2661682	19.12%	1.881%
20	10.369	1213	1221	1239	rBV	2436845	4082487	29.33%	2.885%
21	10.752	1277	1286	1299	rBV	2597754	4431668	31.84%	3.132%
22	10.893	1302	1310	1328	rBV	1983185	3507666	25.20%	2.479%
23	11.663	1436	1441	1442	rBV4	10079	14019	0.10%	0.010%
24	12.157	1521	1525	1531	rVB3	24508	39786	0.29%	0.028%
25	12.334	1550	1555	1560	rBV2	38668	55916	0.40%	0.040%
26	12.457	1570	1576	1582	rBV7	24972	46359	0.33%	0.033%
27	13.104	1683	1686	1692	rVB8	12946	19005	0.14%	0.013%
28	13.346	1717	1727	1730	rBV8	10114	26919	0.19%	0.019%
29	13.393	1730	1735	1741	rVB	48370	75557	0.54%	0.053%
30	13.845	1808	1812	1817	rBV8	9263	15671	0.11%	0.011%
31	13.898	1817	1821	1825	rVV7	15487	25168	0.18%	0.018%
32	13.987	1827	1836	1844	rVV	3699600	5107878	36.70%	3.610%
33	14.051	1844	1847	1852	rVV6	25831	48469	0.35%	0.034%
34	14.098	1852	1855	1860	rVB6	12321	20227	0.15%	0.014%
35	14.275	1878	1885	1894	rVV	4464269	6465785	46.45%	4.570%
36	14.340	1894	1896	1900	rVB4	16917	21775	0.16%	0.015%

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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_P\Methods\SFAM-EPA-BP120722.M
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37	14.469	1914	1918	1921	rVB2	30888	46664	0.34%	0.033%
38	14.581	1931	1937	1945	rBV	3335518	4899241	35.20%	3.463%
39	14.645	1945	1948	1951	rVB5	14301	15406	0.11%	0.011%
40	14.787	1965	1972	1994	rBV	1487658	2531654	18.19%	1.789%

41	14.934	1994	1997	2001	rVV4	30326	49999	0.36%	0.035%
42	14.987	2002	2006	2013	rVB6	40374	61962	0.45%	0.044%
43	15.245	2046	2050	2055	rVB8	13302	20813	0.15%	0.015%
44	15.369	2066	2071	2074	rBV7	13971	28517	0.20%	0.020%
45	15.422	2077	2080	2085	rVB3	18369	23642	0.17%	0.017%

46	15.575	2099	2106	2120	rBV	5475396	7961129	57.20%	5.627%
47	15.698	2121	2127	2138	rBV	1178043	1618153	11.63%	1.144%
48	15.816	2143	2147	2152	rVB4	29958	49100	0.35%	0.035%
49	15.945	2164	2169	2173	rVB	57790	79843	0.57%	0.056%
50	16.051	2183	2187	2190	rVB6	10532	15741	0.11%	0.011%

51	16.151	2200	2204	2210	rVB8	18188	31115	0.22%	0.022%
52	16.287	2223	2227	2229	rBV5	18174	25842	0.19%	0.018%
53	16.510	2261	2265	2272	rVB6	30134	48885	0.35%	0.035%
54	16.722	2298	2301	2308	rBV6	27456	51049	0.37%	0.036%
55	17.122	2367	2369	2373	rBV4	11122	16458	0.12%	0.012%

56	17.339	2400	2406	2417	rBV	4191854	6137846	44.10%	4.338%
57	17.439	2417	2423	2436	rVB	5844970	8444474	60.67%	5.968%
58	17.698	2462	2467	2470	rBV6	32004	57196	0.41%	0.040%
59	17.998	2515	2518	2523	rVB7	17770	23120	0.17%	0.016%
60	18.210	2549	2554	2564	rBV	555232	871359	6.26%	0.616%

61	18.492	2600	2602	2607	rBV5	21446	23999	0.17%	0.017%
62	18.545	2607	2611	2620	rBV2	90399	157628	1.13%	0.111%
63	19.104	2703	2706	2710	rVB	60916	68155	0.49%	0.048%
64	19.251	2727	2731	2736	rVB3	89264	121316	0.87%	0.086%
65	19.322	2738	2743	2744	rBV	104672	141508	1.02%	0.100%

66	19.351	2745	2748	2752	rVB	223610	278100	2.00%	0.197%
67	19.445	2760	2764	2770	rBV2	139046	194110	1.39%	0.137%
68	19.675	2797	2803	2815	rBV	8558394	11669917	83.84%	8.248%
69	19.904	2839	2842	2845	rVB	118296	130481	0.94%	0.092%
70	20.180	2885	2889	2893	rVB	549042	617418	4.44%	0.436%

71	20.298	2905	2909	2913	rVB	682683	751143	5.40%	0.531%
72	20.351	2914	2918	2924	rVV2	148308	198939	1.43%	0.141%
73	20.669	2968	2972	2976	rBV	846780	915084	6.57%	0.647%
74	20.727	2978	2982	2987	rBV	480846	552402	3.97%	0.390%
75	21.122	3045	3049	3060	rBV	2110573	3161267	22.71%	2.234%

76	21.433	3097	3102	3107	rVB	5534786	7225498	51.91%	5.107%
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Title : SVOA CALIBRATION

77	21.569	3121	3125	3135	rVB	936989	1166046	8.38%	0.824%
78	22.039	3201	3205	3207	rBV	990345	1227081	8.82%	0.867%
79	22.069	3207	3210	3220	rVB2	865487	1543736	11.09%	1.091%
80	22.551	3288	3292	3298	rVB	924162	1311231	9.42%	0.927%
81	23.127	3385	3390	3397	rBV2	1640245	2809383	20.18%	1.986%
82	23.757	3489	3497	3512	rVB	6319719	13918527	100.00%	9.837%
83	23.910	3517	3523	3532	rVB2	3760561	7909475	56.83%	5.590%
84	24.527	3623	3628	3637	rVB	1105142	2293376	16.48%	1.621%

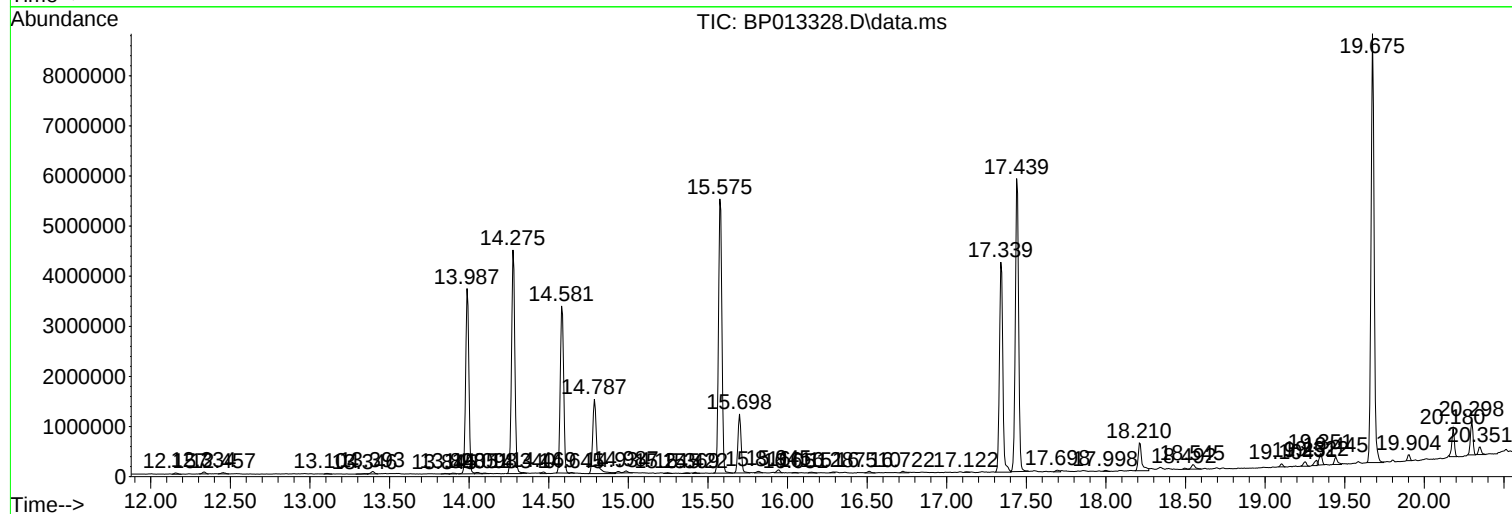
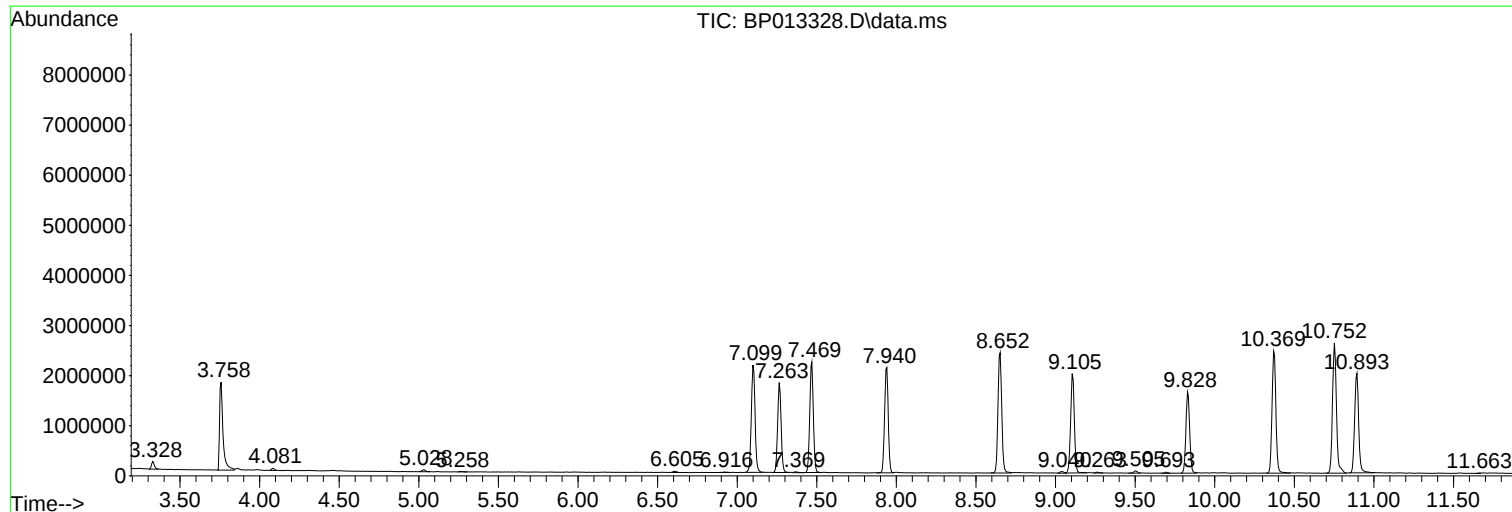
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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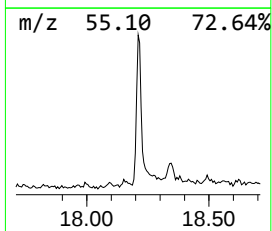
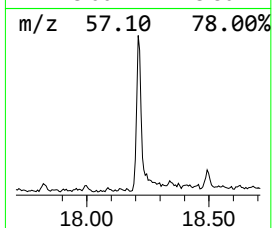
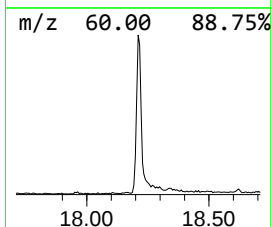
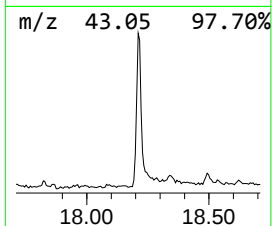
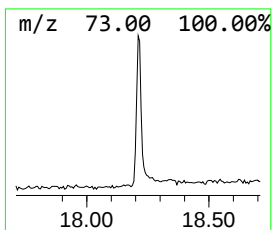
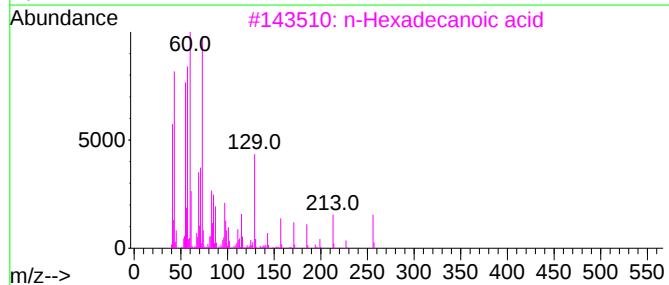
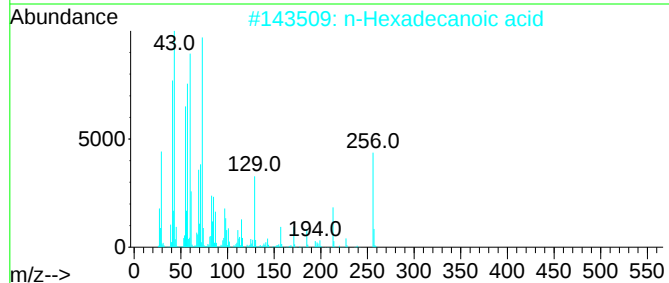
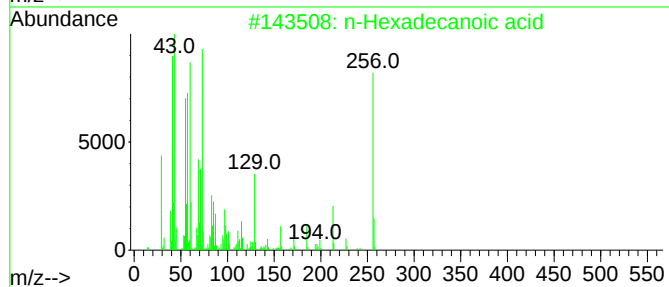
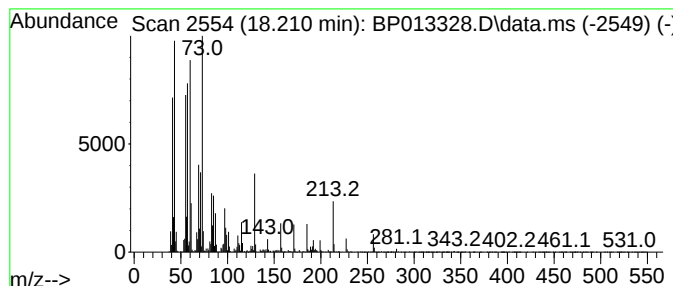
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.210	2.84 ng/ul	871359	Phenanthrene-d10	17.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	93
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91



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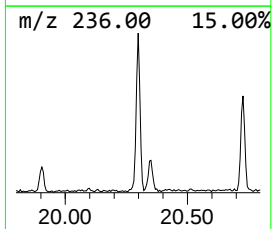
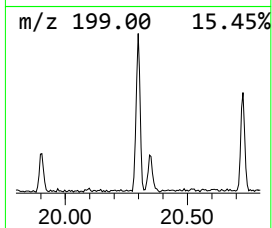
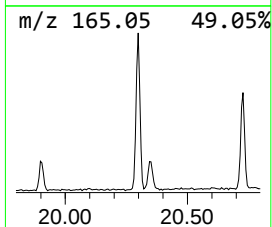
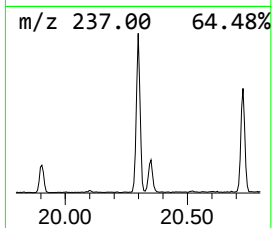
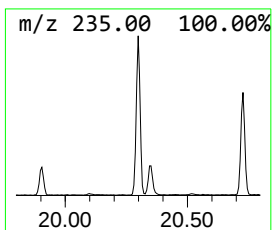
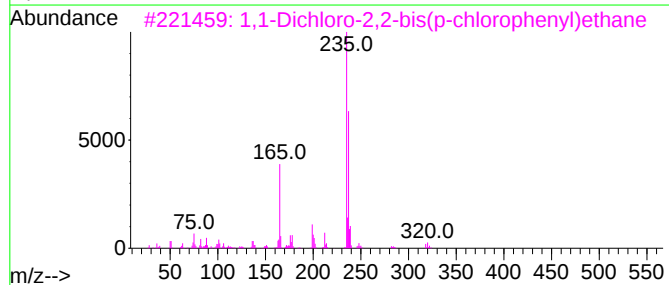
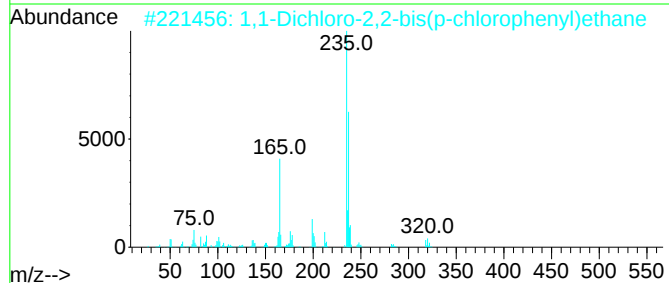
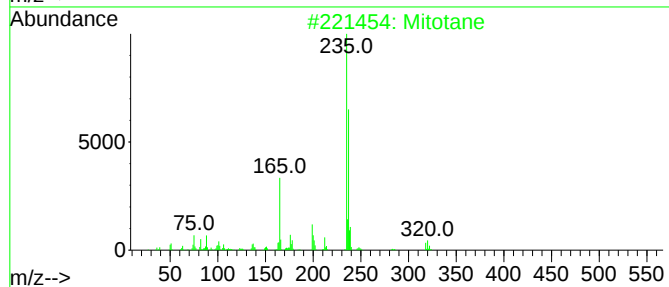
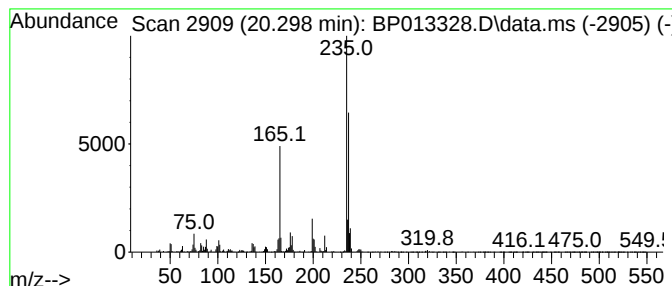
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Mitotane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.298	2.08 ng/ul	751143	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mitotane	318	C14H10Cl4	000053-19-0	99
2			1,1-Dichloro-2,2-bis(p-chlorophe...	318	C14H10Cl4	000072-54-8	99
3			1,1-Dichloro-2,2-bis(p-chlorophe...	318	C14H10Cl4	000072-54-8	91
4			m,p'-DDD	318	C14H10Cl4	004329-12-8	91
5			3-Butanone, 1,1-bis(4-chlorophen...	320	C18H18Cl2O	1000158-20-4	87



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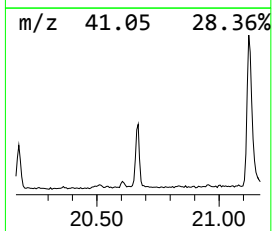
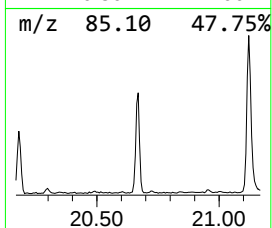
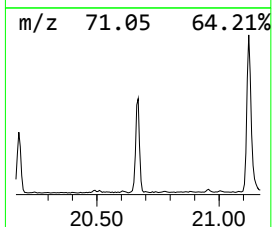
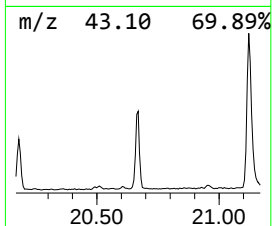
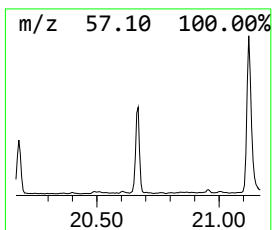
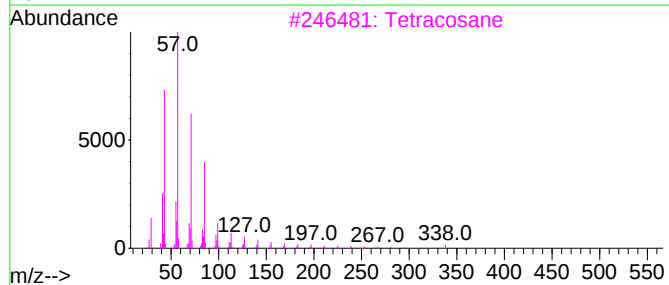
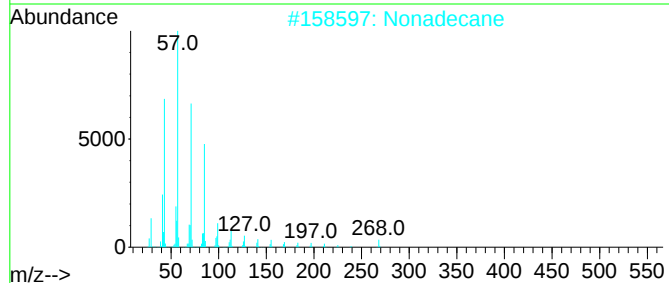
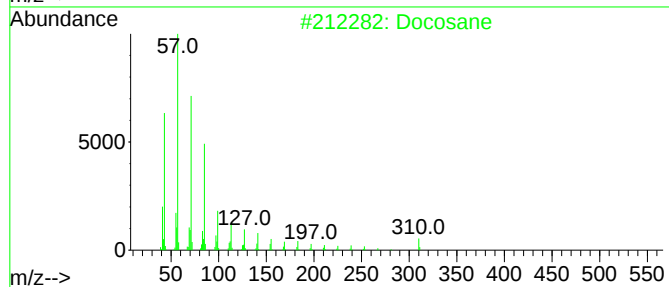
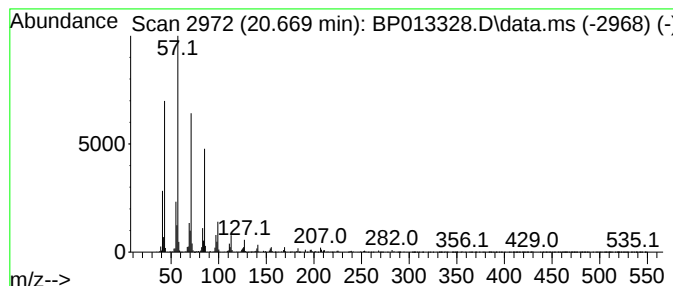
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.669	2.53 ng/ul	915084	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Docosane	310	C22H46	000629-97-0	96
2			Nonadecane	268	C19H40	000629-92-5	95
3			Tetracosane	338	C24H50	000646-31-1	93
4			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	91
5			Octacosane, 1-iodo-	520	C28H57I	1000406-32-2	91



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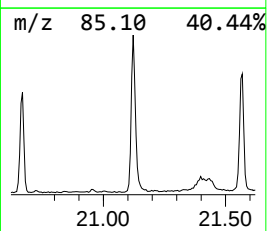
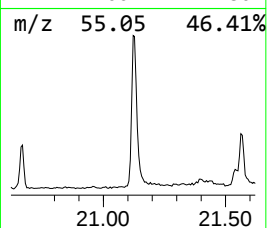
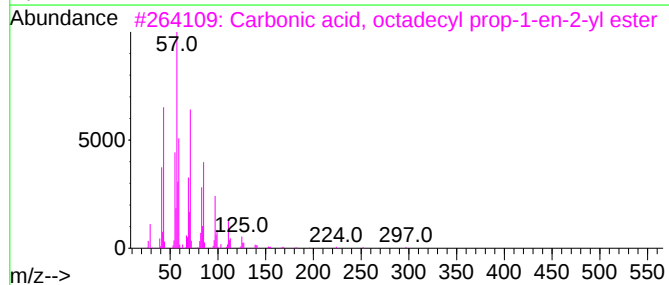
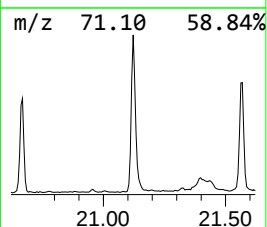
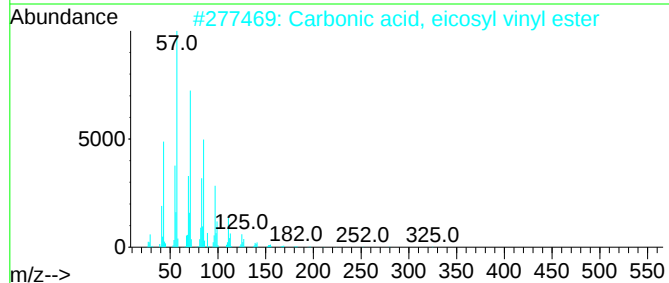
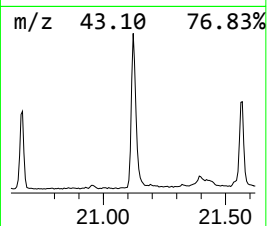
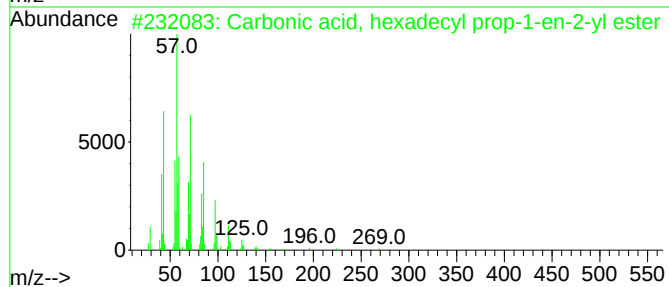
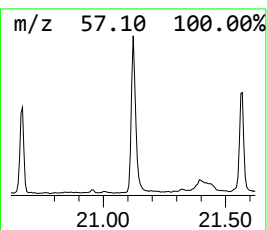
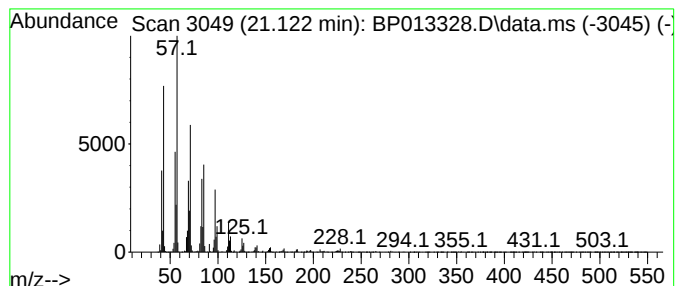
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Carbonic acid, hexadecyl pr... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.122	8.75 ng/ul	3161270	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, hexadecyl prop-1-...	326	C20H38O3	1000382-90-3	91
2			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	91
3			Carbonic acid, octadecyl prop-1-...	354	C22H42O3	1000383-11-5	91
4			Sulfurous acid, octadecyl 2-prop...	376	C21H44O3S	1000309-12-7	87
5			Carbonic acid, octadecyl vinyl e...	340	C21H40O3	1000382-54-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013328.D
Acq On : 25 Dec 2022 08:09
Operator : CG/JU
Sample : N6018-19
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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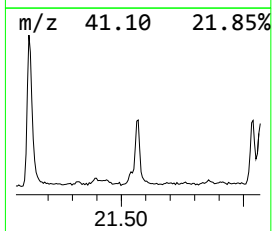
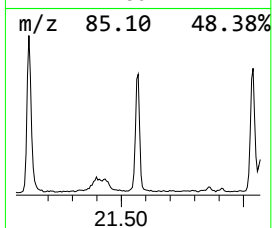
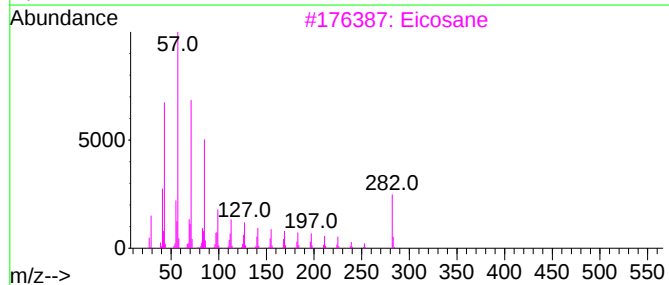
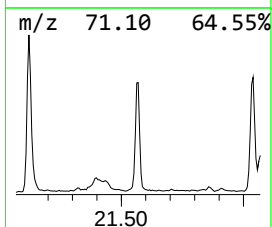
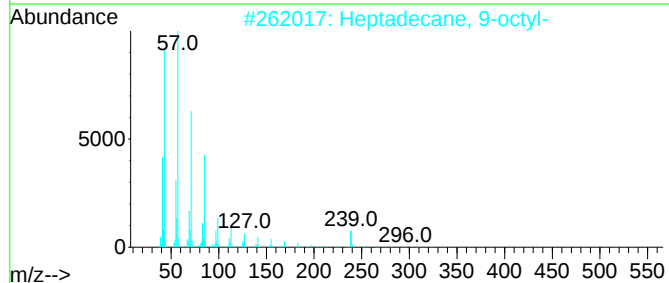
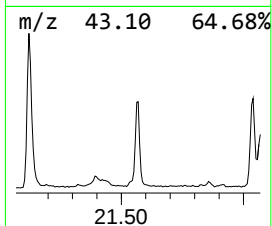
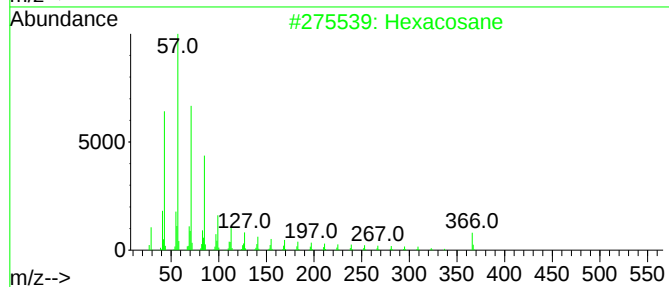
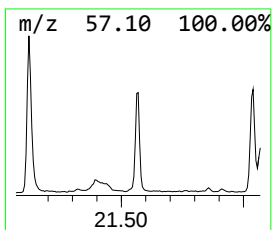
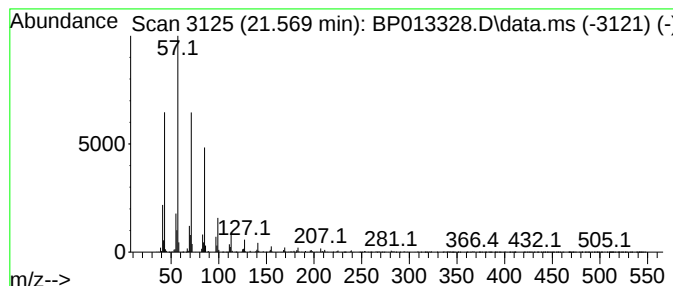
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.569	3.23 ng/ul	1166050	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	94
2			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	93
3			Eicosane	282	C20H42	000112-95-8	90
4			Eicosane, 1-iodo-	408	C20H41I	1000406-31-8	90
5			triacontane	422	C30H62	000638-68-6	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013328.D
Acq On : 25 Dec 2022 08:09
Operator : CG/JU
Sample : N6018-19
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYK4

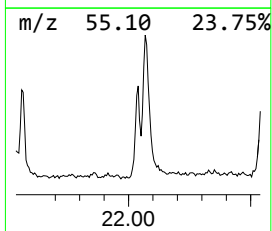
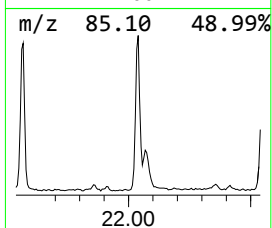
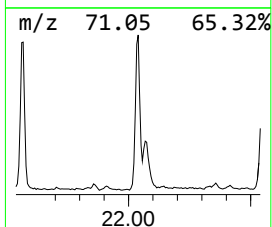
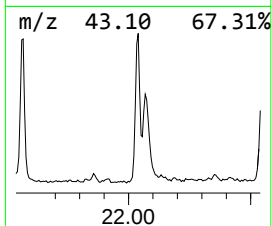
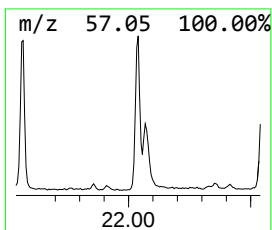
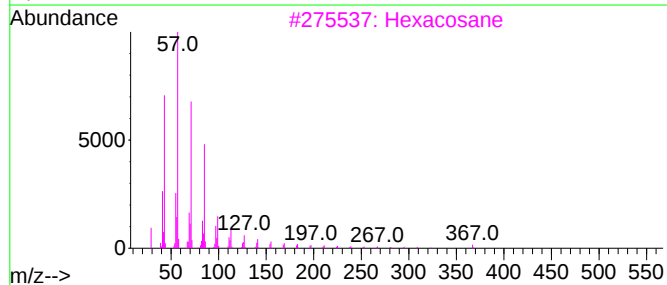
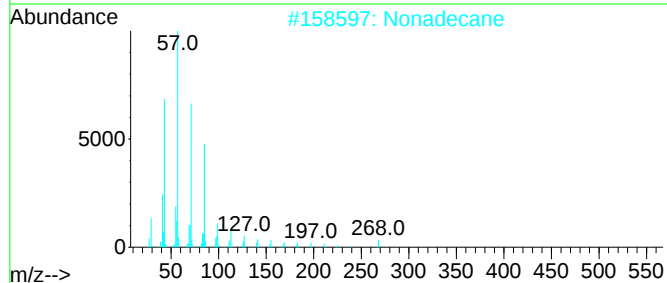
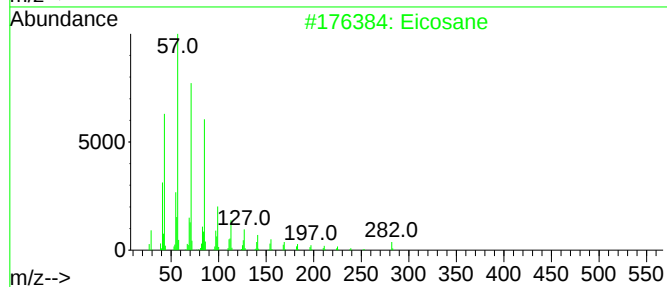
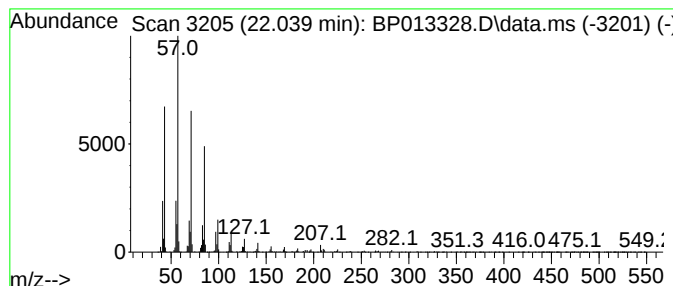
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.039	3.40 ng/ul	1227080	Chrysene-d12	21.433

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane	282	C20H42	000112-95-8	96
2		Nonadecane	268	C19H40	000629-92-5	94
3		Hexacosane	366	C26H54	000630-01-3	94
4		2-Bromo dodecane	248	C12H25Br	013187-99-0	93
5		Docosane, 1-iodo-	436	C22H45I	1000406-31-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013328.D
Acq On : 25 Dec 2022 08:09
Operator : CG/JU
Sample : N6018-19
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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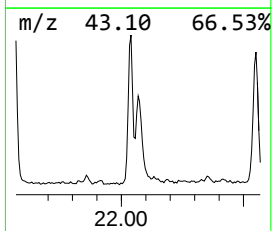
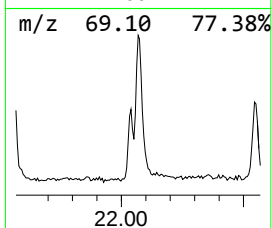
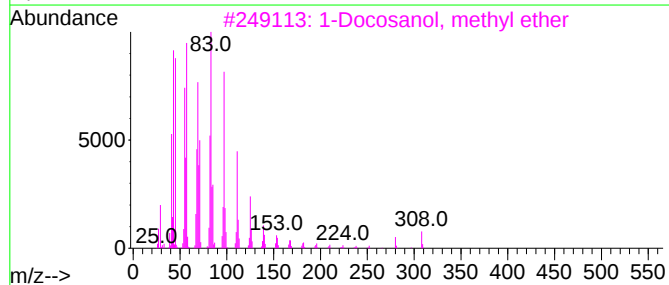
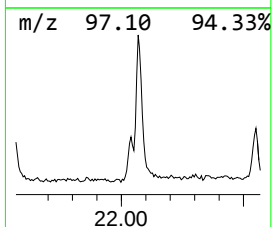
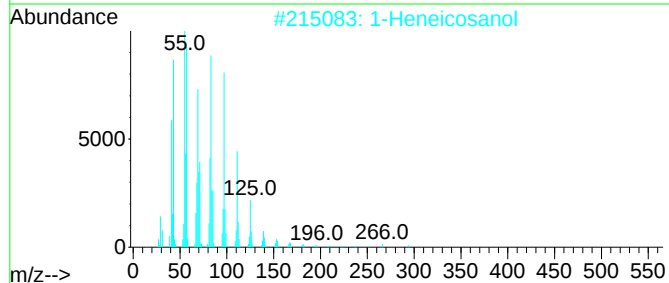
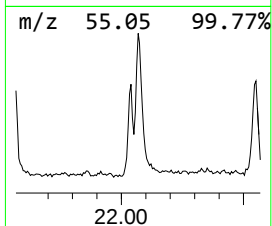
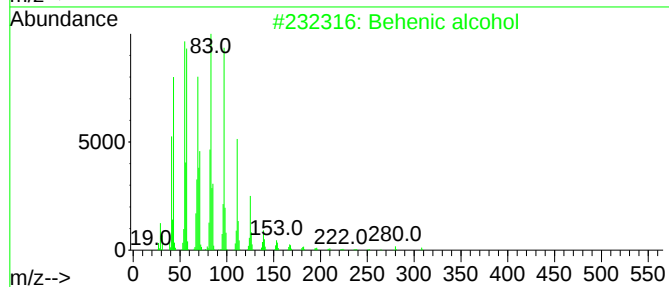
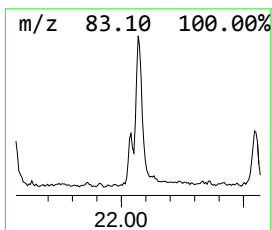
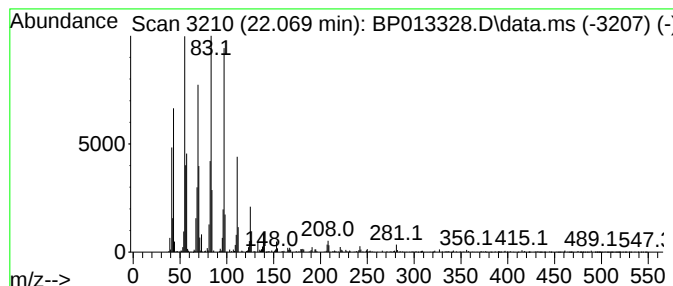
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Behenic alcohol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.069	4.27 ng/ul	1543740	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Behenic alcohol	326	C22H46O	000661-19-8	94
2			1-Heneicosanol	312	C21H44O	015594-90-8	91
3			1-Docosanol, methyl ether	340	C23H48O	1000333-91-9	91
4			n-Nonadecanol-1	284	C19H40O	001454-84-8	90
5			9-Tricosene, (Z)-	322	C23H46	027519-02-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013328.D
Acq On : 25 Dec 2022 08:09
Operator : CG/JU
Sample : N6018-19
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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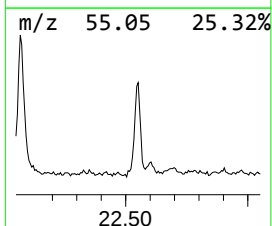
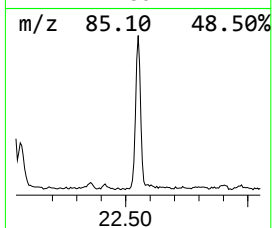
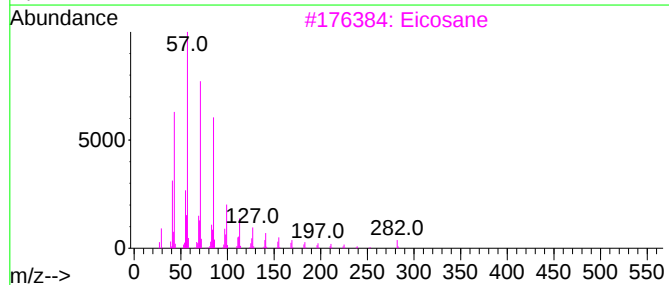
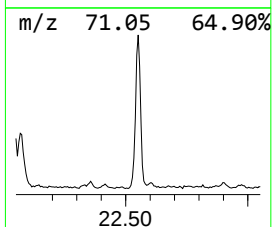
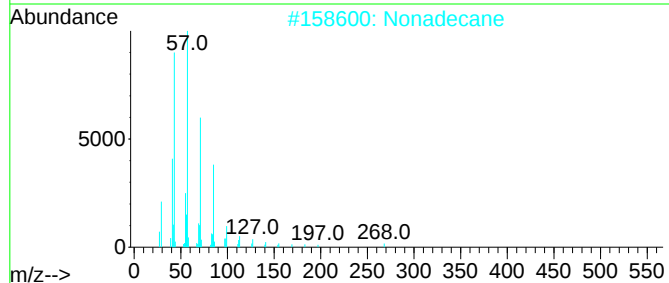
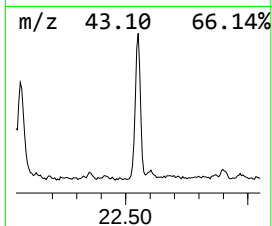
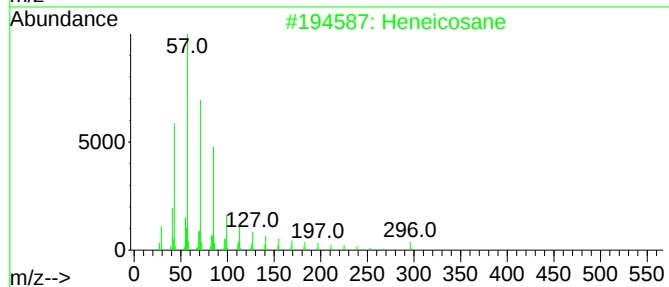
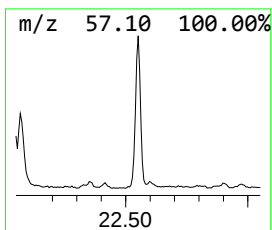
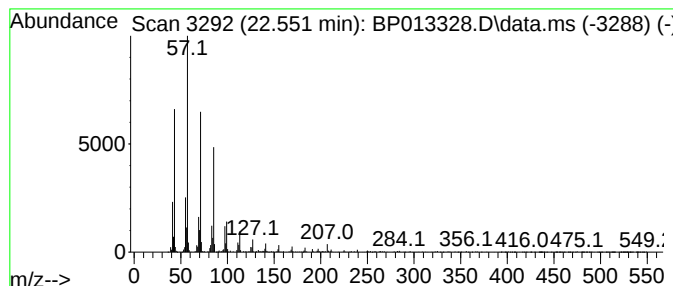
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.551	3.63 ng/ul	1311230	Chrysene-d12	21.433

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heneicosane	296	C21H44	000629-94-7	98
2		Nonadecane	268	C19H40	000629-92-5	96
3		Eicosane	282	C20H42	000112-95-8	95
4		Heneicosane	296	C21H44	000629-94-7	95
5		Eicosane	282	C20H42	000112-95-8	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013328.D
Acq On : 25 Dec 2022 08:09
Operator : CG/JU
Sample : N6018-19
Misc :
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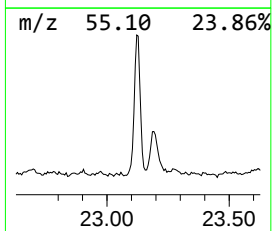
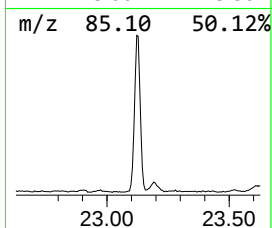
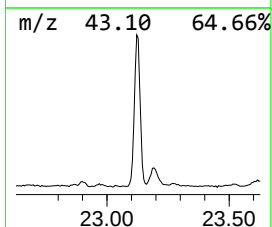
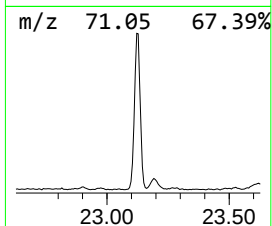
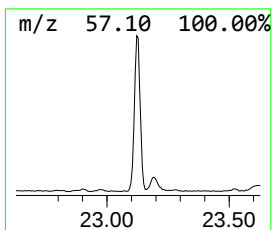
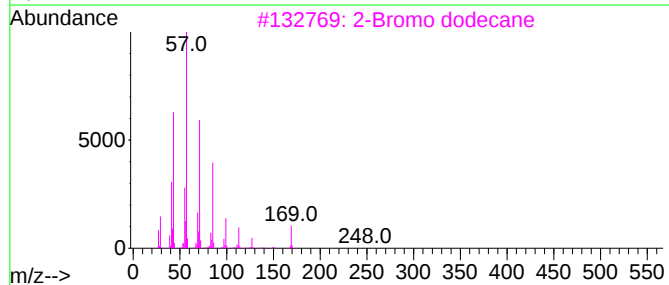
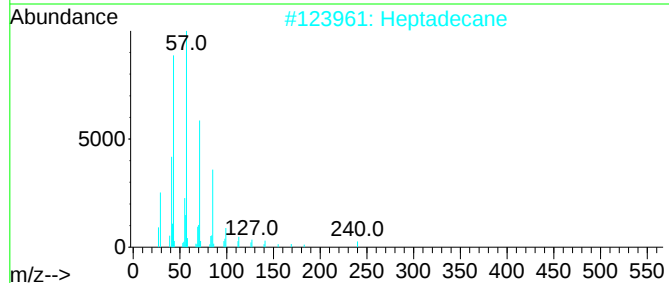
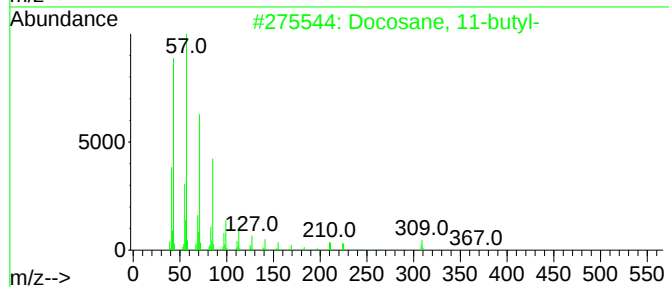
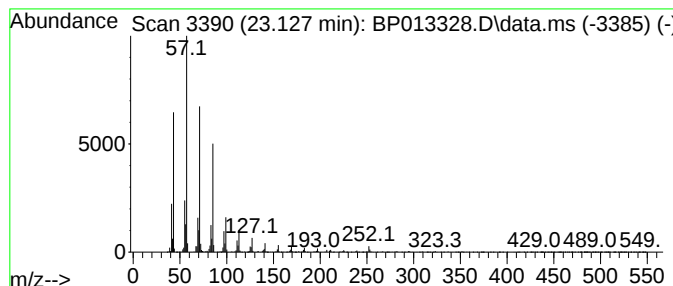
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.127	7.10 ng/ul	2809380	Perylene-d12	23.910	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane, 11-butyl-	366	C26H54	013475-76-8	94
2	Heptadecane	240	C17H36	000629-78-7	93
3	2-Bromo dodecane	248	C12H25Br	013187-99-0	93
4	Eicosane, 3-methyl-	296	C21H44	006418-46-8	91
5	Docosane, 1-iodo-	436	C22H45I	1000406-31-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013328.D
Acq On : 25 Dec 2022 08:09
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Misc :
ALS Vial : 41 Sample Multiplier: 1

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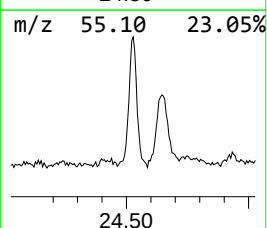
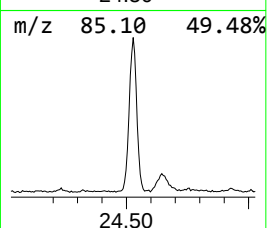
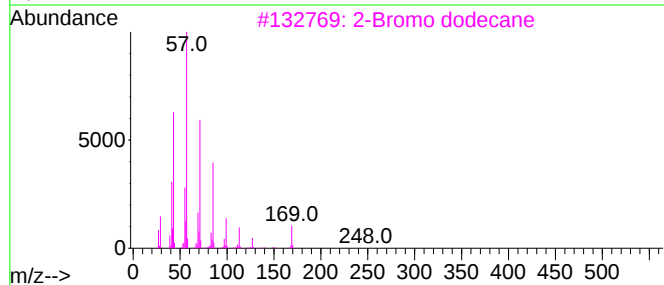
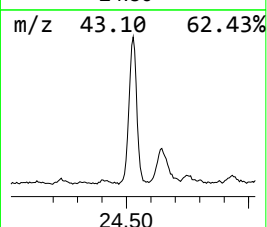
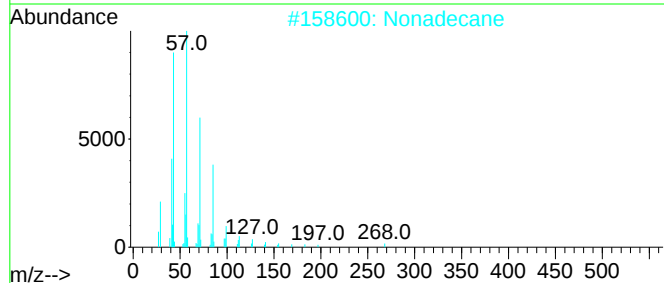
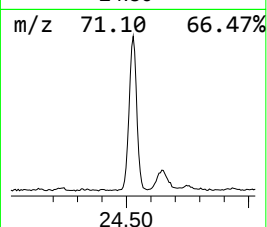
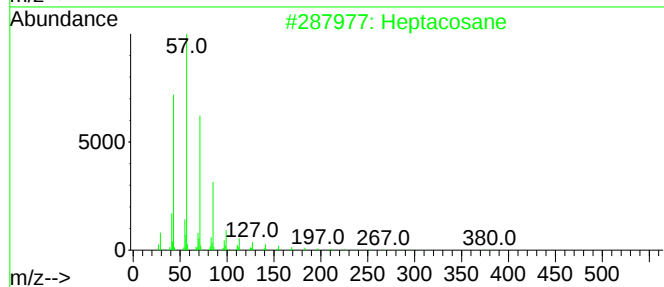
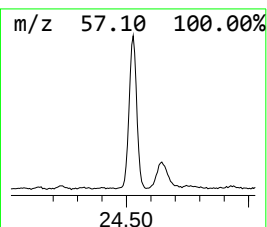
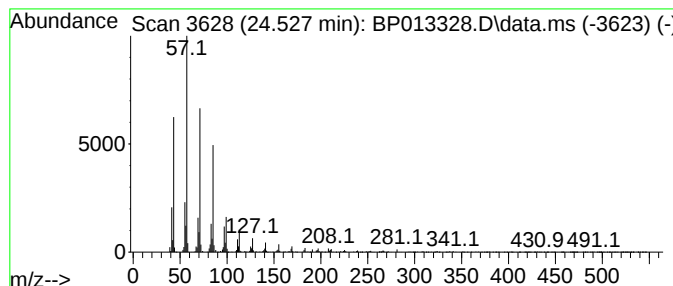
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.527	5.80 ng/ul	2293380	Perylene-d12	23.910

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptacosane	380	C27H56	000593-49-7	96
2		Nonadecane	268	C19H40	000629-92-5	94
3		2-Bromo dodecane	248	C12H25Br	013187-99-0	93
4		Tricosane	324	C23H48	000638-67-5	91
5		Octacosane	394	C28H58	000630-02-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013328.D
Acq On : 25 Dec 2022 08:09
Operator : CG/JU
Sample : N6018-19
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYK4

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.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
n-Hexadecanoic ...	18.210	2.8	ng/ul	871359	4	17.339	6137850	20.0
Mitotane	20.298	2.1	ng/ul	751143	5	21.433	7225500	20.0
(DEL) Alkane: S...	20.669	2.5	ng/ul	915084	5	21.433	7225500	20.0
Carbonic acid, ...	21.122	8.8	ng/ul	3161270	5	21.433	7225500	20.0
(DEL) Alkane: S...	21.569	3.2	ng/ul	1166050	5	21.433	7225500	20.0
(DEL) Alkane: S...	22.039	3.4	ng/ul	1227080	5	21.433	7225500	20.0
Behenic alcohol	22.069	4.3	ng/ul	1543740	5	21.433	7225500	20.0
(DEL) Alkane: S...	22.551	3.6	ng/ul	1311230	5	21.433	7225500	20.0
(DEL) Alkane: S...	23.127	7.1	ng/ul	2809380	6	23.910	7909480	20.0
(DEL) Alkane: S...	24.527	5.8	ng/ul	2293380	6	23.910	7909480	20.0