

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122022\
 Data File : BM038147.D
 Acq On : 20 Dec 2022 22:53
 Operator : CG/JU
 Sample : N6008-17
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBXX9

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-BM120722.M
 Title : SVOA CALIBRATION

Signal : TIC: BM038147.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.399	32	36	48	rVB	30774	56690	1.21%	0.099%
2	3.716	86	90	93	rBV5	4157	7334	0.16%	0.013%
3	3.834	106	110	126	rBV	394584	695320	14.86%	1.214%
4	4.157	162	165	172	rVB2	8978	15525	0.33%	0.027%
5	5.122	323	329	336	rVB3	11471	18970	0.41%	0.033%
6	5.363	367	370	377	rBV6	3486	6622	0.14%	0.012%
7	7.028	649	653	656	rBV4	6625	10722	0.23%	0.019%
8	7.204	676	683	696	rBV	664958	1072719	22.93%	1.873%
9	7.369	705	711	722	rBV	510202	771397	16.49%	1.347%
10	7.569	738	745	753	rBV	654762	1037995	22.19%	1.812%
11	8.045	818	826	833	rBV	630109	990893	21.18%	1.730%
12	8.757	938	947	956	rBV	874051	1409897	30.14%	2.461%
13	9.216	1017	1025	1037	rBV	693096	1088031	23.26%	1.899%
14	9.316	1037	1042	1046	rBV8	4957	7662	0.16%	0.013%
15	9.369	1046	1051	1055	rVB6	5879	10553	0.23%	0.018%
16	9.598	1086	1090	1094	rBV3	6935	10162	0.22%	0.018%
17	9.939	1140	1148	1160	rBV	580492	952781	20.36%	1.663%
18	10.157	1180	1185	1192	rVB7	4632	9547	0.20%	0.017%
19	10.422	1226	1230	1233	rBV5	4232	7121	0.15%	0.012%
20	10.480	1233	1240	1251	rVB	983630	1617243	34.57%	2.823%
21	10.857	1295	1304	1313	rBV	915130	1519238	32.47%	2.652%
22	10.998	1320	1328	1346	rBV	814109	1388417	29.68%	2.424%
23	11.392	1391	1395	1399	rVB6	3429	5399	0.12%	0.009%
24	11.639	1432	1437	1447	rVB5	9716	16912	0.36%	0.030%
25	12.251	1538	1541	1547	rVB6	5359	8788	0.19%	0.015%
26	12.439	1568	1573	1578	rBV2	16795	25813	0.55%	0.045%
27	12.945	1653	1659	1662	rBV8	2829	4704	0.10%	0.008%
28	13.274	1711	1715	1718	rBV5	4424	6639	0.14%	0.012%
29	13.480	1745	1750	1754	rBV5	7786	11434	0.24%	0.020%
30	13.663	1777	1781	1785	rVB2	10757	13505	0.29%	0.024%
31	14.080	1843	1852	1865	rVV	1862085	2453976	52.45%	4.284%
32	14.192	1867	1871	1875	rVB6	3783	5678	0.12%	0.010%
33	14.304	1886	1890	1894	rBV7	5730	9975	0.21%	0.017%
34	14.368	1894	1901	1908	rBV	2179883	3040850	65.00%	5.309%
35	14.545	1925	1931	1935	rBV7	6877	11296	0.24%	0.020%
36	14.674	1946	1953	1960	rBV	1429654	2054048	43.90%	3.586%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120722.M
 Title : SVOA CALIBRATION

37	14.768	1966	1969	1973	rVB6	3697	5721	0.12%	0.010%
38	14.868	1979	1986	1998	rBV	801403	1159243	24.78%	2.024%
39	14.957	1998	2001	2005	rVV2	19894	25955	0.55%	0.045%
40	15.021	2009	2012	2016	rVV4	9344	15110	0.32%	0.026%

41	15.068	2017	2020	2029	rVB8	18088	31275	0.67%	0.055%
42	15.445	2078	2084	2087	rBV7	3488	6846	0.15%	0.012%
43	15.486	2087	2091	2096	rVB7	4155	8320	0.18%	0.015%
44	15.610	2108	2112	2113	rBV4	4001	4945	0.11%	0.009%
45	15.657	2114	2120	2128	rVB	2731762	3782810	80.85%	6.604%

46	15.780	2135	2141	2149	rBV	734272	997813	21.33%	1.742%
47	15.892	2157	2160	2165	rVB3	14164	19664	0.42%	0.034%
48	15.980	2172	2175	2177	rBV4	4522	4973	0.11%	0.009%
49	16.021	2177	2182	2189	rVB	82127	111392	2.38%	0.194%
50	16.192	2208	2211	2213	rBV4	5878	7120	0.15%	0.012%

51	16.280	2224	2226	2231	rVB5	4812	5264	0.11%	0.009%
52	16.351	2236	2238	2240	rBV3	5521	6335	0.14%	0.011%
53	16.509	2260	2265	2268	rBV7	7482	13210	0.28%	0.023%
54	16.580	2273	2277	2280	rBV4	8706	11499	0.25%	0.020%
55	16.798	2310	2314	2318	rBV4	19221	27108	0.58%	0.047%

56	16.904	2330	2332	2337	rBV6	3705	6846	0.15%	0.012%
57	17.145	2369	2373	2375	rBV4	8414	11272	0.24%	0.020%
58	17.292	2394	2398	2405	rBV6	25542	52107	1.11%	0.091%
59	17.356	2405	2409	2412	rVV4	14983	20867	0.45%	0.036%
60	17.409	2412	2418	2427	rVV	1793295	2500515	53.45%	4.365%

61	17.509	2427	2435	2441	rVV	2930115	3988384	85.25%	6.963%
62	17.562	2442	2444	2447	rVV4	19802	24678	0.53%	0.043%
63	17.604	2447	2451	2455	rVB7	17407	31299	0.67%	0.055%
64	17.680	2461	2464	2468	rBV4	9170	9279	0.20%	0.016%
65	17.751	2473	2476	2479	rBV5	23269	25107	0.54%	0.044%

66	17.880	2495	2498	2502	rBV6	6626	10053	0.21%	0.018%
67	17.927	2502	2506	2510	rVB6	12683	17494	0.37%	0.031%
68	18.056	2526	2528	2535	rVB7	14418	19178	0.41%	0.033%
69	18.168	2542	2547	2553	rBV3	38666	72559	1.55%	0.127%
70	18.227	2553	2557	2562	rBV	47731	80689	1.72%	0.141%

71	18.286	2562	2567	2577	rBV	819550	1127380	24.10%	1.968%
72	18.633	2622	2626	2632	rVB4	20449	36379	0.78%	0.064%
73	18.703	2635	2638	2644	rBV5	28626	39041	0.83%	0.068%
74	18.868	2663	2666	2672	rVB7	24872	36934	0.79%	0.064%
75	18.945	2677	2679	2683	rBV5	13331	17147	0.37%	0.030%

76	19.203	2719	2723	2728	rBV5	23685	43608	0.93%	0.076%
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77	19.292	2735	2738	2743	rBV4	20813	31172	0.67%	0.054%
78	19.356	2745	2749	2754	rBV2	40699	55823	1.19%	0.097%
79	19.415	2755	2759	2760	rBV2	96236	123022	2.63%	0.215%
80	19.550	2778	2782	2791	rBV	266922	388267	8.30%	0.678%
81	19.792	2817	2823	2833	rBV	3707416	4678545	100.00%	8.168%
82	19.939	2844	2848	2856	rBV	149639	190930	4.08%	0.333%
83	20.303	2903	2910	2915	rBV4	65782	155968	3.33%	0.272%
84	20.633	2963	2966	2967	rBV2	47713	54907	1.17%	0.096%
85	20.756	2984	2987	2992	rVB	166170	195407	4.18%	0.341%
86	21.262	3069	3073	3081	rVB	758154	969311	20.72%	1.692%
87	21.468	3104	3108	3110	rBV	254098	318332	6.80%	0.556%
88	21.574	3122	3126	3132	rVB	1947396	2421018	51.75%	4.226%
89	23.233	3405	3408	3414	rVB	188817	267583	5.72%	0.467%
90	23.868	3509	3516	3526	rVB2	1939258	3808754	81.41%	6.649%
91	24.027	3537	3543	3554	rVB2	1018948	2077106	44.40%	3.626%
92	24.609	3638	3642	3651	rVB	330220	621675	13.29%	1.085%
93	28.538	4302	4310	4332	rVB5	906349	3328550	71.14%	5.811%
94	29.044	4390	4396	4421	rVB10	648592	2806429	59.99%	4.899%

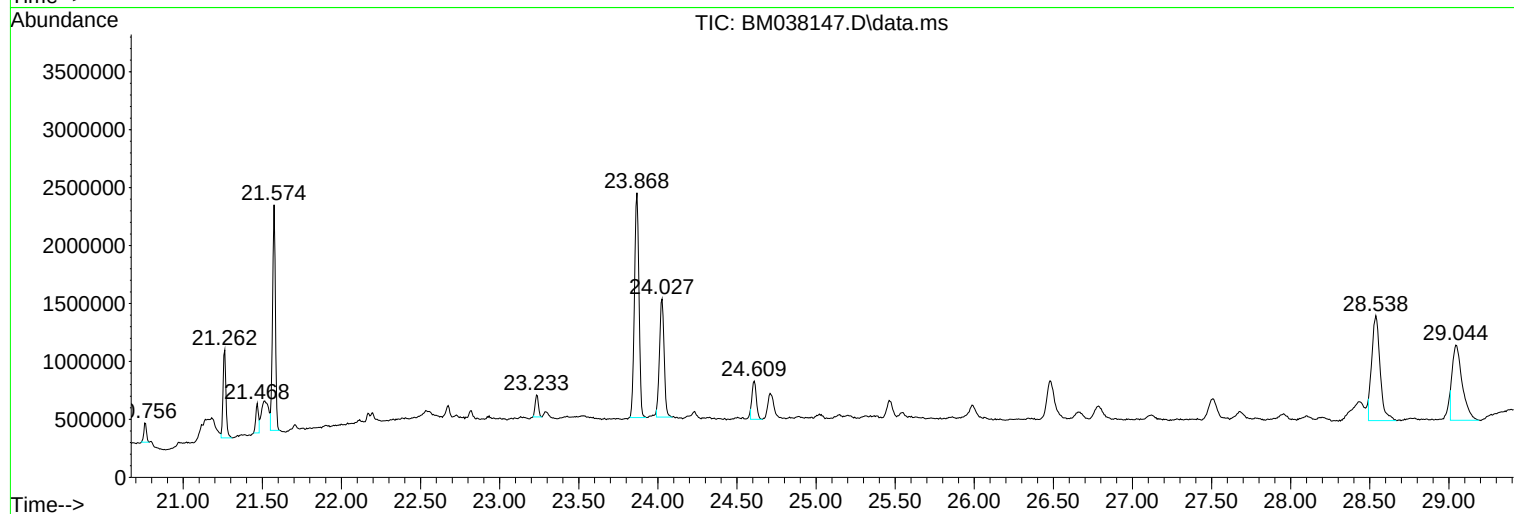
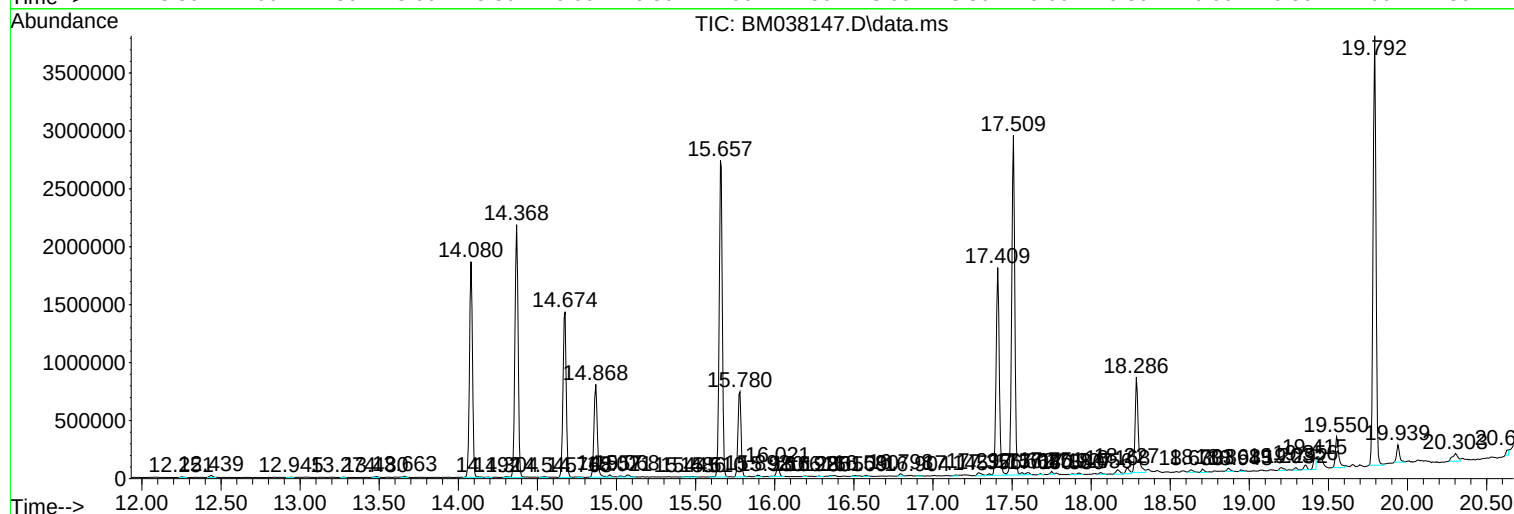
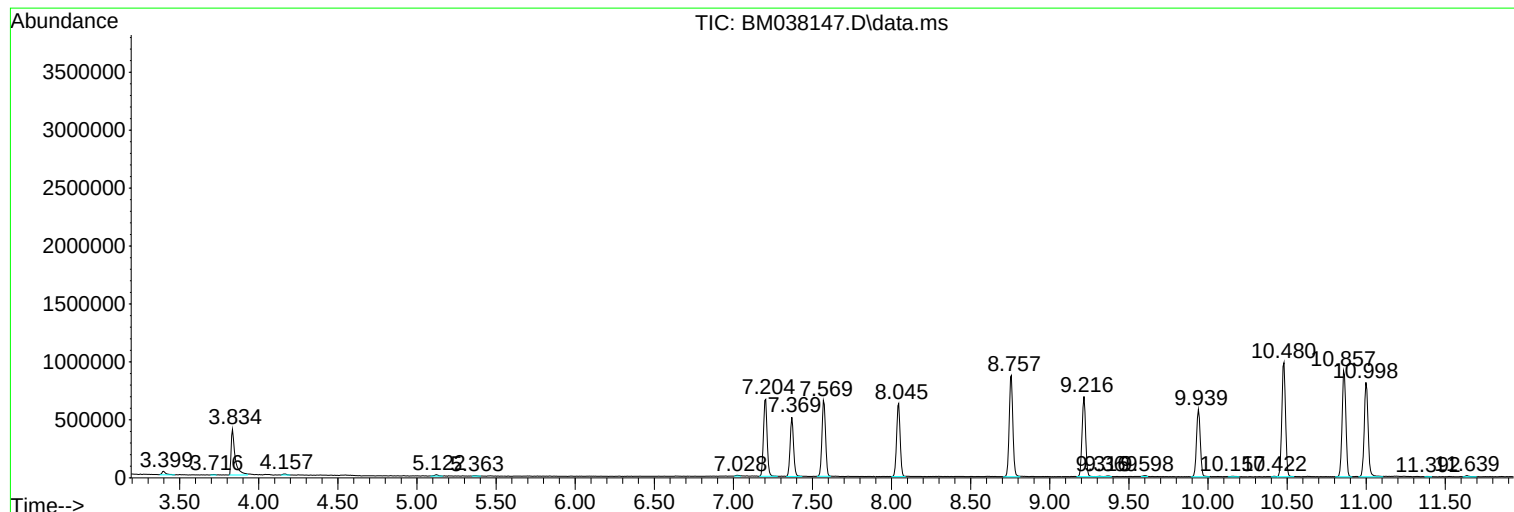
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120722.M
Quant Title : SVOA CALIBRATION

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



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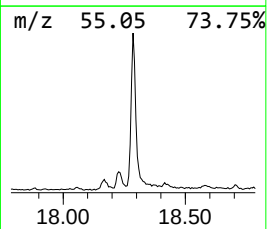
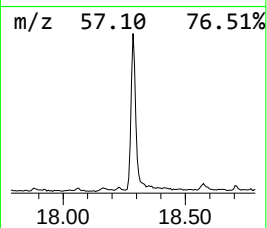
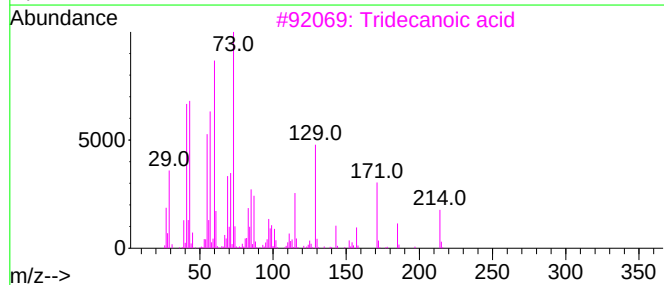
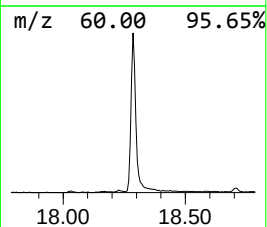
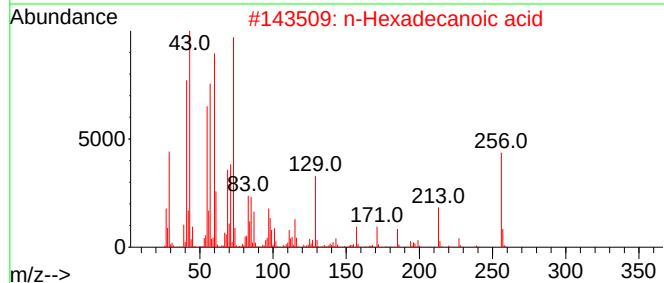
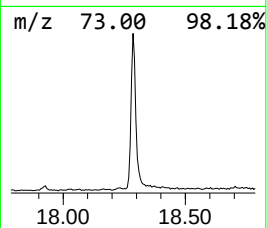
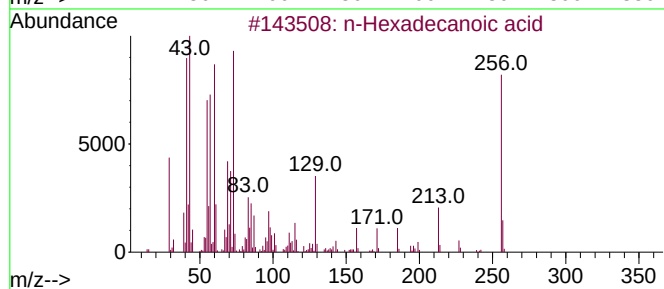
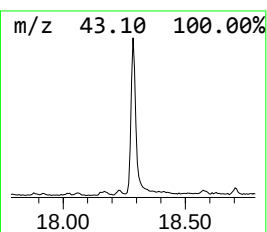
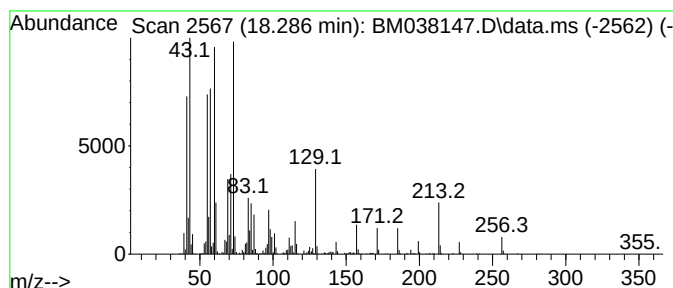
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.286	9.02 ng/ul	1127380	Phenanthrene-d10	17.409

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			Tridecanoic acid	214	C13H26O2	000638-53-9	96
4			Tridecanoic acid	214	C13H26O2	000638-53-9	94
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94



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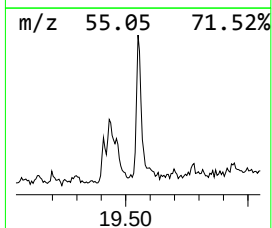
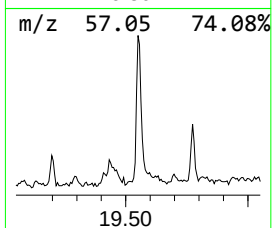
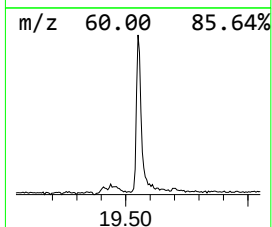
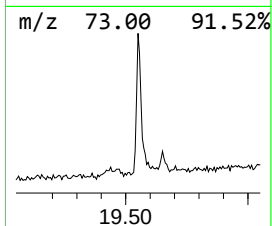
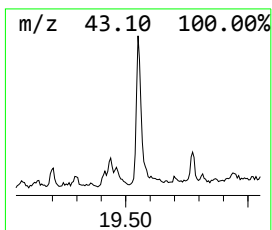
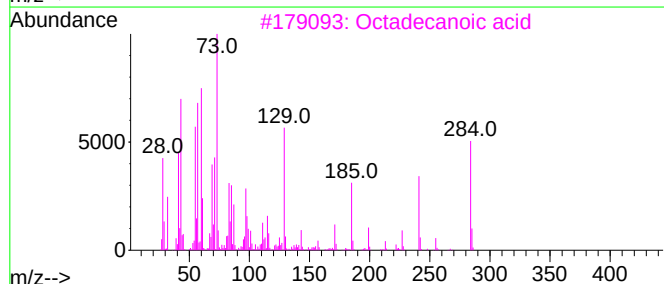
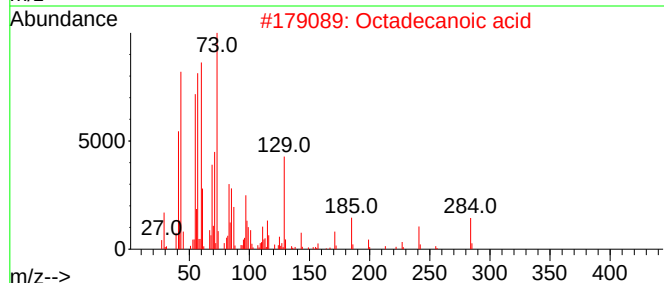
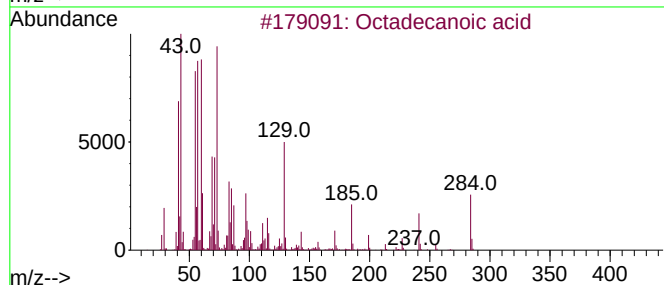
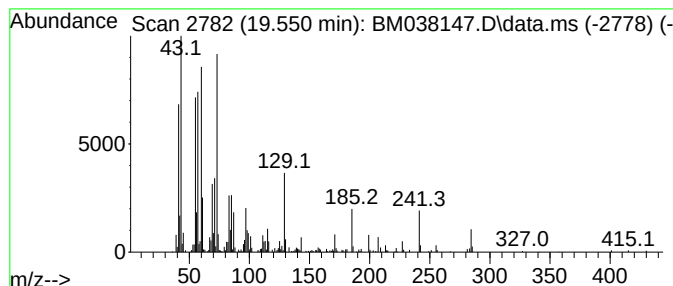
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.551	3.21 ng/ul	388267	Chrysene-d12	21.574

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	99
4			Octadecanoic acid	284	C18H36O2	000057-11-4	98
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	81



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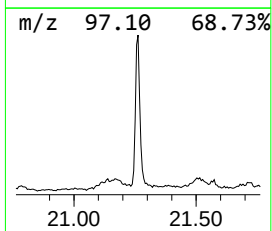
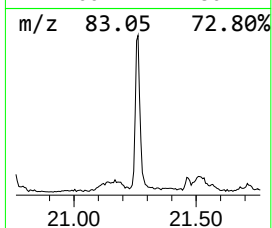
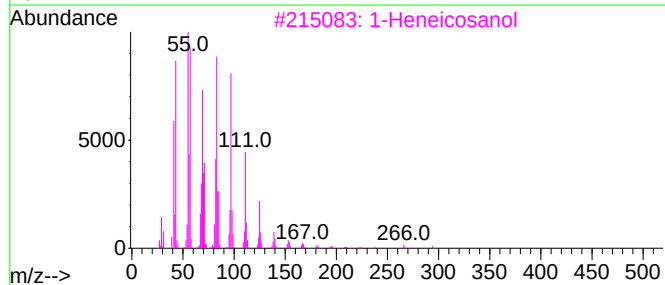
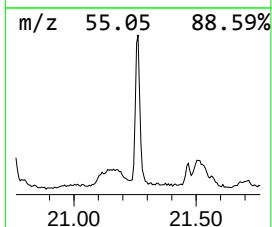
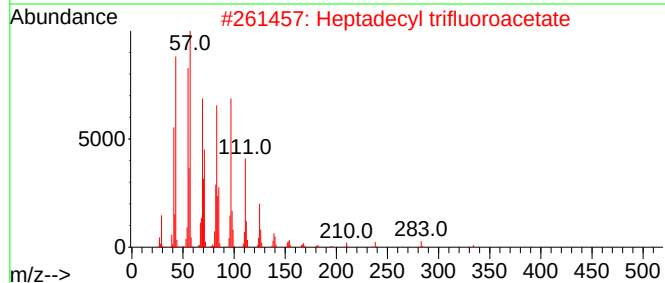
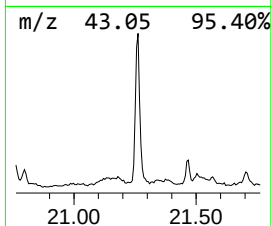
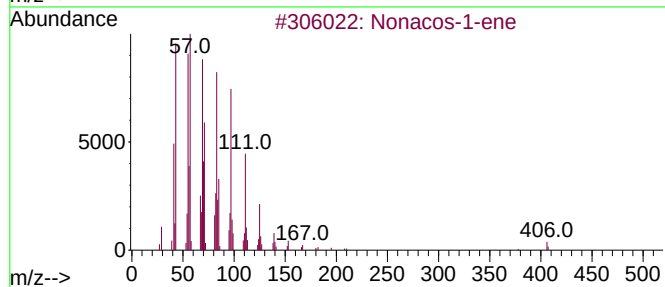
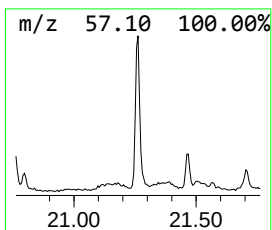
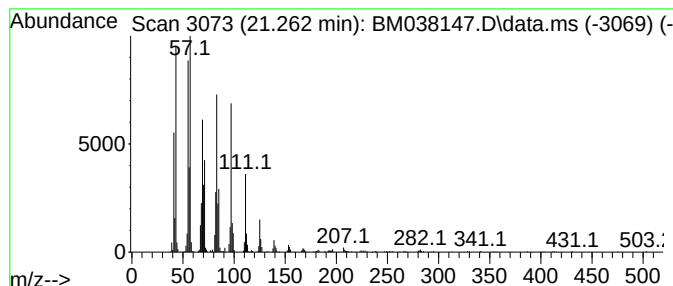
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.262	8.01 ng/ul	969311	Chrysene-d12	21.574	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonacos-1-ene	406	C29H58	018835-35-3	94
2	Heptadecyl trifluoroacetate	352	C19H35F3O2	1010351-87-0	94
3	1-Heneicosanol	312	C21H44O	015594-90-8	91
4	Pentacos-1-ene	350	C25H50	016980-85-1	91
5	Octacosanol	410	C28H58O	000557-61-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122022\
Data File : BM038147.D
Acq On : 20 Dec 2022 22:53
Operator : CG/JU
Sample : N6008-17
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DBXX9

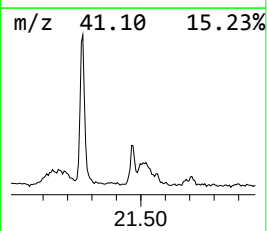
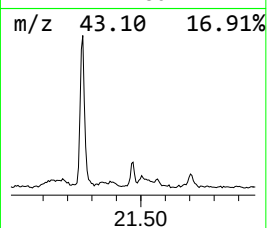
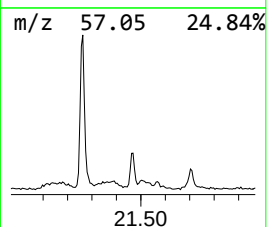
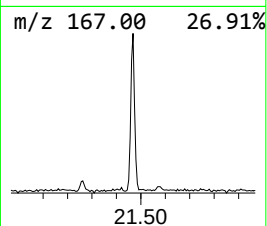
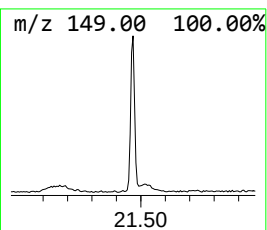
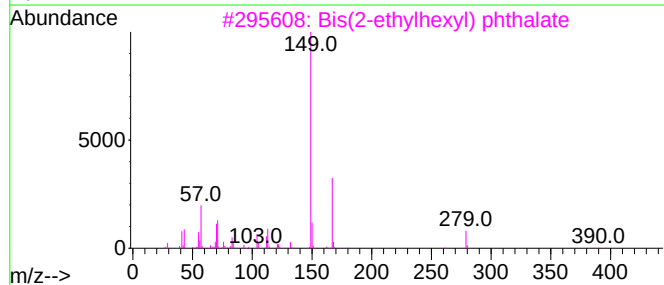
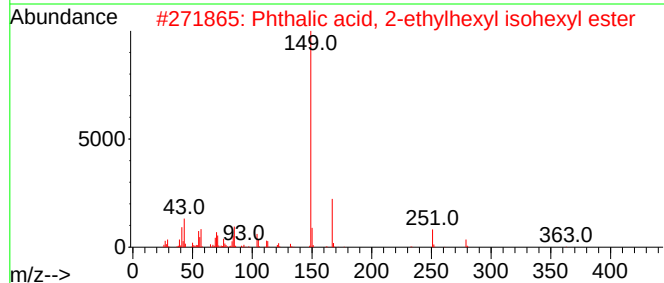
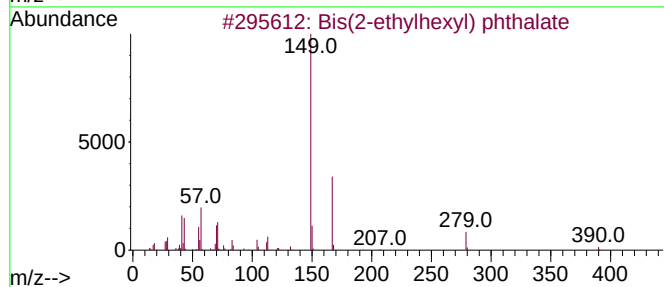
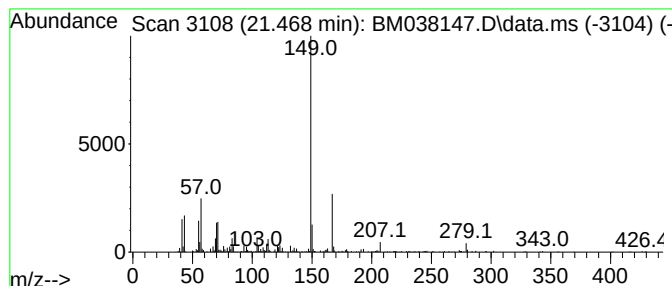
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Bis(2-ethylhexyl) phthalate Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.468	2.63 ng/ul	318332	Chrysene-d12	21.574

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	87
2			Phthalic acid, 2-ethylhexyl isoh...	362	C22H34O4	1000308-98-5	86
3			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	80
4			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	80
5			Phthalic acid, 2-ethylhexyl tetr...	474	C30H50O4	1000309-00-4	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122022\
Data File : BM038147.D
Acq On : 20 Dec 2022 22:53
Operator : CG/JU
Sample : N6008-17
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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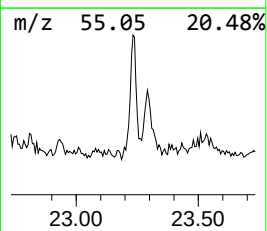
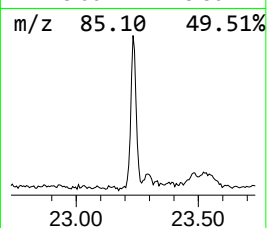
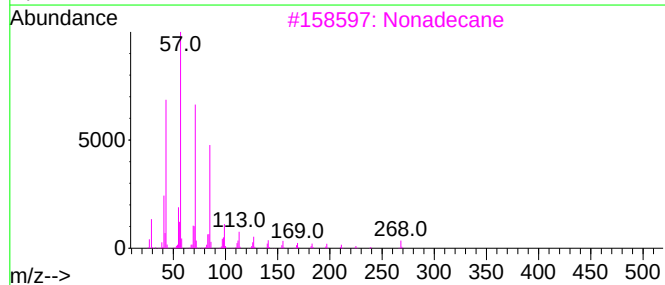
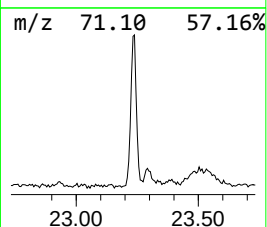
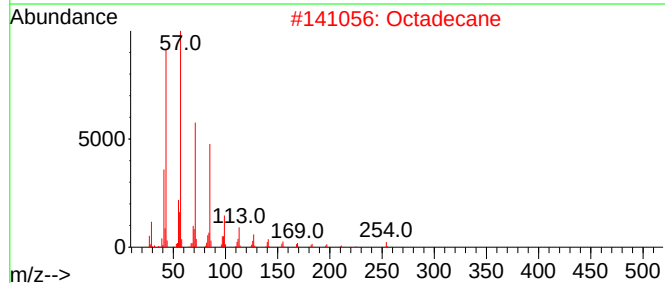
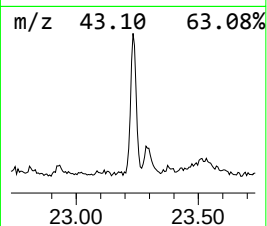
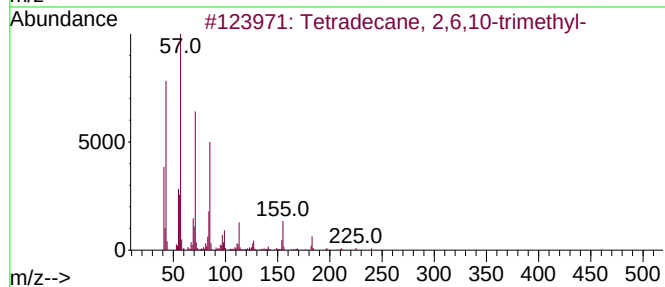
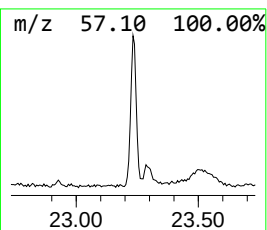
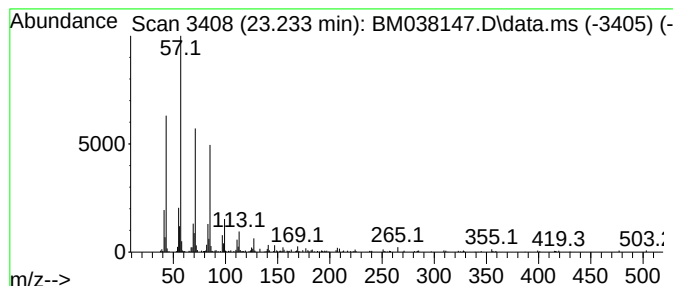
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.233	2.58 ng/ul	267583	Perylene-d12	24.027

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	91
2		Octadecane	254	C18H38	000593-45-3	87
3		Nonadecane	268	C19H40	000629-92-5	86
4		Tetracosane	338	C24H50	000646-31-1	86
5		Tridecane, 1-iodo-	310	C13H27I	035599-77-0	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122022\
Data File : BM038147.D
Acq On : 20 Dec 2022 22:53
Operator : CG/JU
Sample : N6008-17
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DBXX9

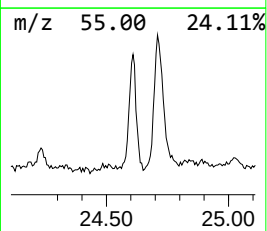
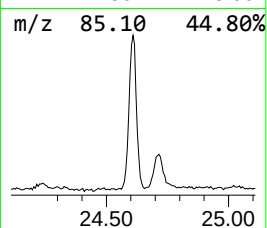
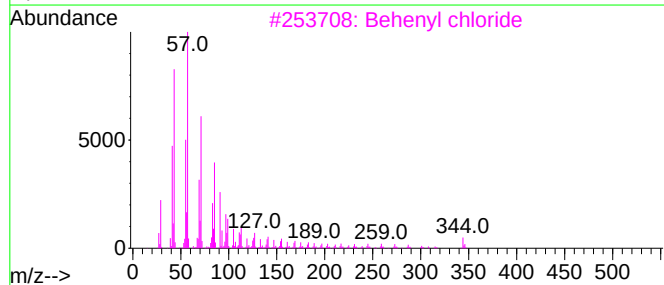
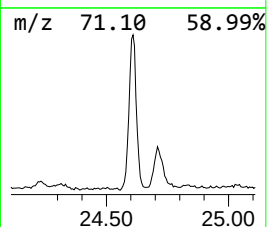
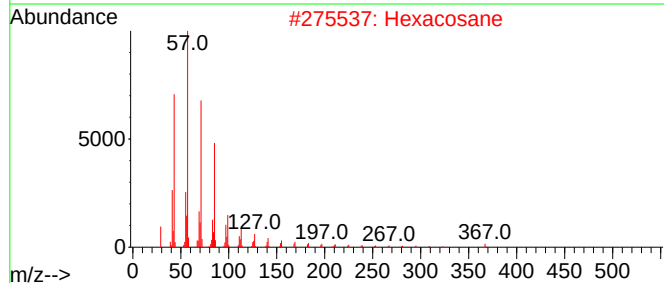
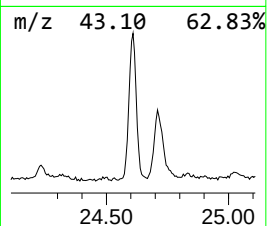
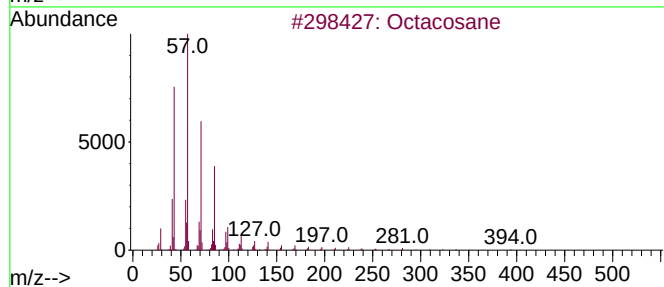
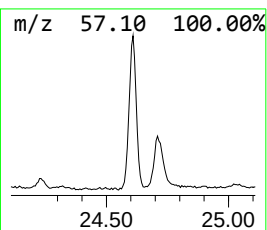
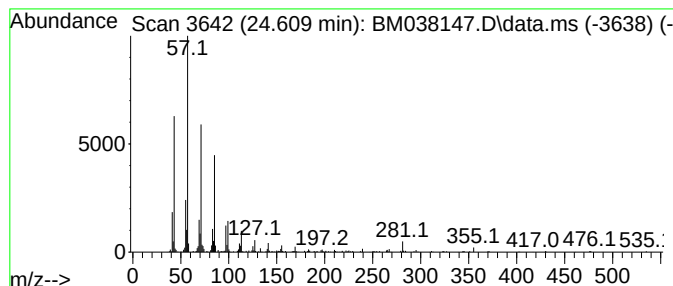
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.609	5.99 ng/ul	621675	Perylene-d12	24.027

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octacosane	394	C28H58	000630-02-4	90
2		Hexacosane	366	C26H54	000630-01-3	87
3		Behenyl chloride	344	C22H45Cl	042217-03-8	87
4		Tricosane	324	C23H48	000638-67-5	87
5		Tricosane	324	C23H48	000638-67-5	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122022\
Data File : BM038147.D
Acq On : 20 Dec 2022 22:53
Operator : CG/JU
Sample : N6008-17
Misc :
ALS Vial : 14 Sample Multiplier: 1

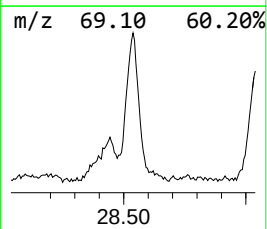
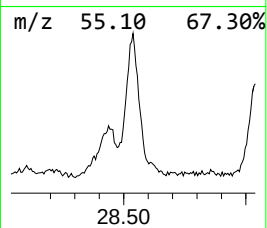
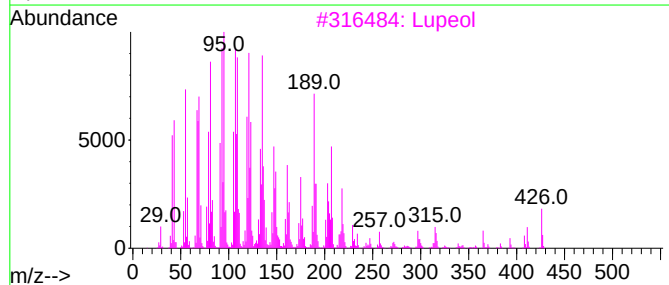
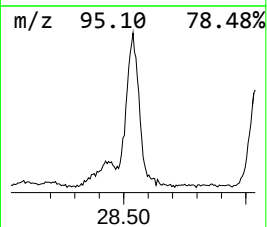
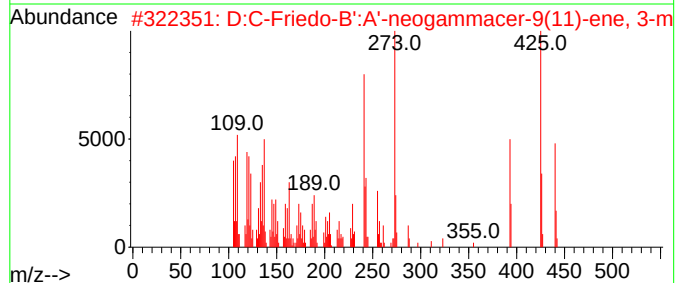
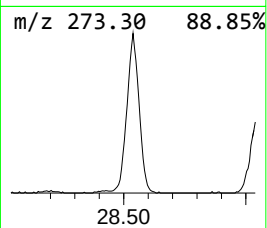
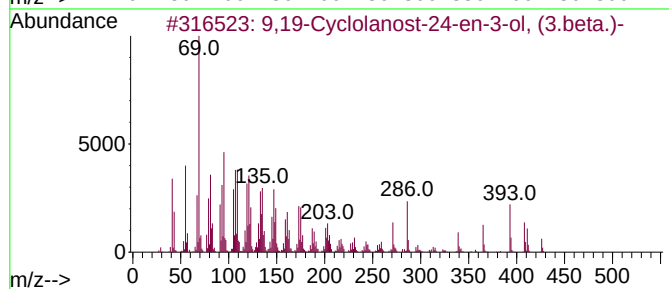
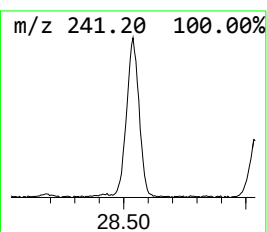
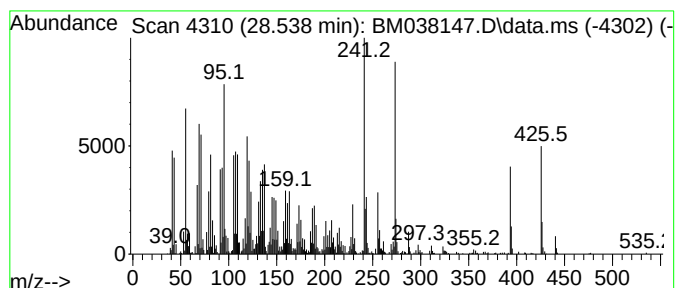
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 9,19-Cyclolanost-24-en-3-ol... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
28.538	32.05 ng/ul	3328550	Perylene-d12	24.027	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,19-Cyclolanost-24-en-3-ol, (3...	426	C30H50O	000469-38-5	56
2	D:C-Friedo-B':A'-neogammacer-9(1...	440	C31H52O	004555-56-0	47
3	Lupeol	426	C30H50O	000545-47-1	35
4	Pregnan-18-ol, (5.alpha.)-	304	C21H36O	056053-09-9	22
5	9,19-Cyclolanost-25-en-3-ol, 24-...	440	C31H52O	000511-61-5	20



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122022\
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Acq On : 20 Dec 2022 22:53
Operator : CG/JU
Sample : N6008-17
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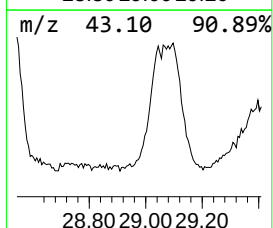
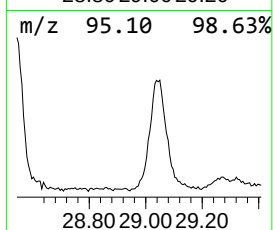
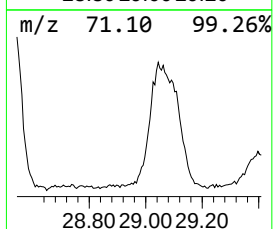
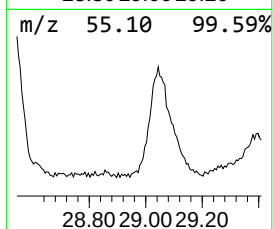
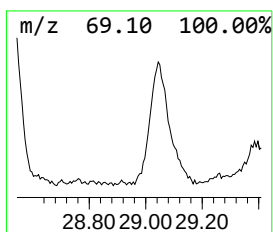
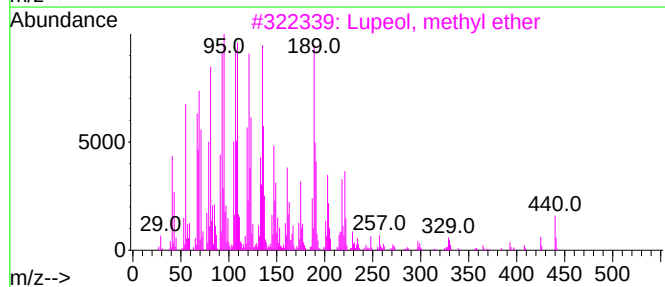
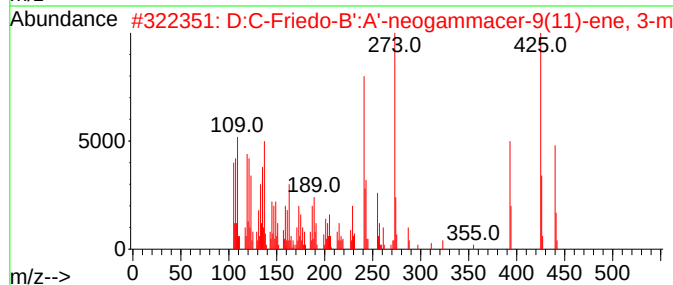
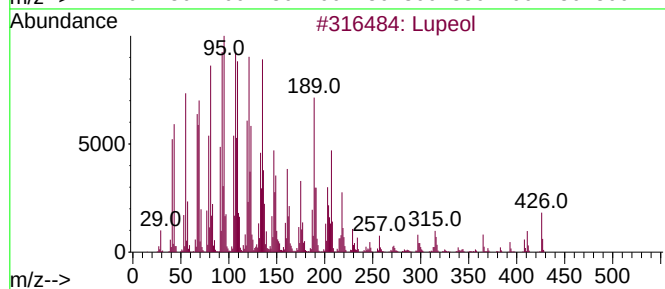
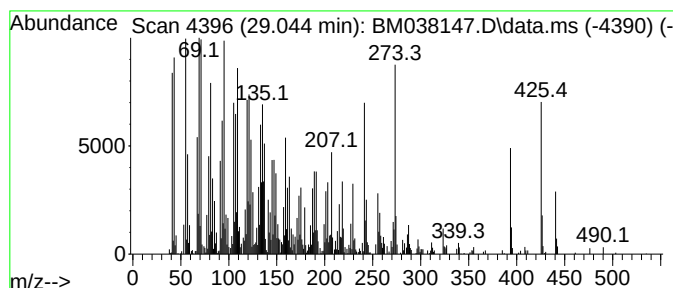
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 Lupeol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.044	27.02 ng/ul	2806430	Perylene-d12	24.027
Hit#	of 5	Tentative ID	MW MolForm	CAS# Qual
1		Lupeol	426 C30H50O	000545-47-1 53
2		D:C-Friedo-B':A'-neogammacer-9(1...	440 C31H52O	004555-56-0 46
3		Lupeol, methyl ether	440 C31H52O	025279-38-3 38
4		1,2-Dihydro-4H-2-(p-anisyl)-3,1-...	241 C15H15NO2	082085-87-8 25
5		Crinan-1-ol, (1.alpha.)-	273 C16H19NO3	041928-89-6 18



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122022\
Data File : BM038147.D
Acq On : 20 Dec 2022 22:53
Operator : CG/JU
Sample : N6008-17
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120722.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 ...	18.286	9.0	ng/ul	1127380	4	17.409	2500520	20.0
Octadecanoic acid	19.551	3.2	ng/ul	388267	5	21.574	2421020	20.0
(DEL) Alkane: S...	21.262	8.0	ng/ul	969311	5	21.574	2421020	20.0
Bis(2-ethylhexy...	21.468	2.6	ng/ul	318332	5	21.574	2421020	20.0
(DEL) Alkane: S...	23.233	2.6	ng/ul	267583	6	24.027	2077110	20.0
(DEL) Alkane: S...	24.609	6.0	ng/ul	621675	6	24.027	2077110	20.0
9,19-Cyclolanos...	28.538	32.0	ng/ul	3328550	6	24.027	2077110	20.0
Lupeol	29.044	27.0	ng/ul	2806430	6	24.027	2077110	20.0