

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
 Data File : BP019024.D
 Acq On : 22 Dec 2023 05:48
 Operator : MA/JU
 Sample : 05808-19
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0B15

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-BP120123.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP019024.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.252	25	28	32	rVB2	16217	18985	0.19%	0.026%
2	3.534	73	76	87	rVB	40704	58886	0.58%	0.082%
3	3.681	98	101	114	rBV	44952	71294	0.70%	0.099%
4	3.781	114	118	123	rVB	17044	19885	0.19%	0.028%
5	3.899	134	138	145	rVB	16575	26640	0.26%	0.037%
6	3.993	151	154	160	rVB	9609	13827	0.14%	0.019%
7	4.917	307	311	319	rVB	27424	44201	0.43%	0.062%
8	5.805	456	462	469	rBV	28669	47055	0.46%	0.065%
9	6.963	653	659	660	rBV	76016	94068	0.92%	0.131%
10	6.993	660	664	668	rVV	203777	314011	3.07%	0.437%
11	7.158	686	692	702	rVB	212476	331863	3.24%	0.462%
12	7.352	718	725	734	rVV	274824	429147	4.20%	0.597%
13	7.440	734	740	748	rVB2	16076	33180	0.32%	0.046%
14	7.816	797	804	815	rBV	553108	868160	8.49%	1.208%
15	8.534	913	926	937	rBV	266968	445451	4.35%	0.620%
16	8.646	937	945	956	rVV	968794	1540282	15.06%	2.144%
17	8.763	959	965	981	rVB	79268	161325	1.58%	0.225%
18	8.981	995	1002	1009	rBV	251901	415858	4.07%	0.579%
19	9.093	1016	1021	1026	rVB3	13163	24594	0.24%	0.034%
20	9.222	1037	1043	1048	rVB3	8432	13961	0.14%	0.019%
21	9.552	1094	1099	1103	rBV2	15661	23034	0.23%	0.032%
22	9.593	1103	1106	1113	rVB6	7370	12205	0.12%	0.017%
23	9.699	1117	1124	1131	rBV	205364	338288	3.31%	0.471%
24	9.916	1146	1161	1175	rBV	249920	712552	6.97%	0.992%
25	10.063	1179	1186	1197	rVB	128811	220493	2.16%	0.307%
26	10.175	1200	1205	1209	rBV2	10322	16505	0.16%	0.023%
27	10.240	1209	1216	1225	rVV	343236	571638	5.59%	0.796%
28	10.475	1253	1256	1265	rVB10	7117	13262	0.13%	0.018%
29	10.610	1271	1279	1288	rBV	713799	1190742	11.64%	1.657%
30	10.763	1299	1305	1320	rBV	21716	56932	0.56%	0.079%
31	11.169	1365	1374	1392	rBV3	133400	423241	4.14%	0.589%
32	11.363	1402	1407	1410	rBV2	35592	62574	0.61%	0.087%
33	11.404	1410	1414	1430	rVB	85311	164649	1.61%	0.229%
34	11.657	1452	1457	1463	rVB	18571	31339	0.31%	0.044%
35	11.981	1506	1512	1518	rBV	39087	68997	0.67%	0.096%
36	12.198	1540	1549	1556	rVB8	11177	29924	0.29%	0.042%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
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37	12.340	1569	1573	1581	rVB8	8168	15498	0.15%	0.022%
38	13.251	1723	1728	1730	rBV6	10733	15895	0.16%	0.022%
39	13.351	1741	1745	1749	rBV2	17622	30078	0.29%	0.042%
40	13.745	1805	1812	1814	rBV8	7790	14989	0.15%	0.021%
41	13.775	1814	1817	1823	rVV8	8744	16594	0.16%	0.023%
42	13.869	1827	1833	1841	rVV	640790	878215	8.59%	1.222%
43	14.145	1871	1880	1891	rVV	698672	1054230	10.31%	1.467%
44	14.322	1900	1910	1919	rBV3	18469	46755	0.46%	0.065%
45	14.451	1924	1932	1941	rBV	1155208	1699714	16.62%	2.366%
46	14.675	1962	1970	1990	rBV	229161	463665	4.53%	0.645%
47	14.822	1990	1995	2000	rVV	25326	41955	0.41%	0.058%
48	14.875	2000	2004	2008	rVV5	16030	23898	0.23%	0.033%
49	15.269	2065	2071	2075	rBV9	5261	12012	0.12%	0.017%
50	15.445	2095	2101	2112	rVB	963149	1398313	13.67%	1.946%
51	15.575	2117	2123	2132	rBV	250924	357172	3.49%	0.497%
52	16.169	2217	2224	2229	rBV9	7711	17786	0.17%	0.025%
53	16.610	2294	2299	2304	rBV5	23095	34084	0.33%	0.047%
54	16.669	2305	2309	2313	rVV	22508	29629	0.29%	0.041%
55	17.010	2363	2367	2370	rVV5	8367	13498	0.13%	0.019%
56	17.204	2394	2400	2406	rBV	1535647	2139733	20.92%	2.978%
57	17.304	2411	2417	2427	rVV	974926	1413713	13.82%	1.968%
58	17.381	2427	2430	2437	rVB8	9107	15555	0.15%	0.022%
59	17.510	2450	2452	2460	rVB3	8643	12869	0.13%	0.018%
60	17.604	2465	2468	2475	rVB9	13743	25296	0.25%	0.035%
61	17.704	2481	2485	2496	rBV	60980	83610	0.82%	0.116%
62	17.881	2511	2515	2521	rBV5	21157	36236	0.35%	0.050%
63	17.975	2526	2531	2538	rVB	72013	94279	0.92%	0.131%
64	18.039	2538	2542	2547	rBV5	22404	34934	0.34%	0.049%
65	18.098	2547	2552	2562	rBV	478410	679229	6.64%	0.945%
66	18.328	2588	2591	2597	rBV	40097	47946	0.47%	0.067%
67	18.386	2597	2601	2604	rBV3	25458	35434	0.35%	0.049%
68	18.510	2618	2622	2625	rBV6	21591	28502	0.28%	0.040%
69	18.616	2636	2640	2642	rBV	48469	60565	0.59%	0.084%
70	18.651	2642	2646	2652	rVB	385404	473494	4.63%	0.659%
71	18.869	2679	2683	2687	rBV	353057	431238	4.22%	0.600%
72	18.916	2687	2691	2697	rVV	149277	219938	2.15%	0.306%
73	18.992	2701	2704	2706	rBV2	42508	54898	0.54%	0.076%
74	19.122	2712	2726	2740	rVB4	334490	1790207	17.50%	2.492%
75	19.333	2758	2762	2773	rBV	227052	324457	3.17%	0.452%
76	19.545	2792	2798	2805	rBV	1478189	1955432	19.12%	2.722%

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77	19.716	2824	2827	2833	rVB	50166	73816	0.72%	0.103%
78	19.845	2846	2849	2854	rVB2	69454	89173	0.87%	0.124%
79	20.051	2880	2884	2893	rBV3	221971	440838	4.31%	0.614%
80	20.386	2938	2941	2943	rBV3	54282	70139	0.69%	0.098%
81	20.733	2996	3000	3002	rBV3	156334	230978	2.26%	0.322%
82	20.833	3009	3017	3019	rBV2	521034	1153047	11.27%	1.605%
83	21.016	3043	3048	3057	rBV	1814394	2566947	25.10%	3.573%
84	21.298	3091	3096	3111	rBV	2148073	3167877	30.97%	4.410%
85	21.633	3150	3153	3157	rVB	230791	252349	2.47%	0.351%
86	21.922	3194	3202	3212	rBV	3631159	6522120	63.76%	9.079%
87	22.251	3254	3258	3267	rBV3	468647	787338	7.70%	1.096%
88	22.433	3285	3289	3296	rVB3	301781	536274	5.24%	0.746%
89	22.633	3319	3323	3328	rVB	606971	842991	8.24%	1.173%
90	22.821	3352	3355	3361	rVB	278408	410571	4.01%	0.572%
91	22.933	3369	3374	3379	rBV	1034385	1523688	14.90%	2.121%
92	22.992	3379	3384	3393	rVB	794427	1470767	14.38%	2.047%
93	23.504	3465	3471	3477	rBV	888420	1779012	17.39%	2.476%
94	23.657	3491	3497	3504	rVB	1380339	2748367	26.87%	3.826%
95	23.892	3532	3537	3544	rVB2	870818	1635014	15.98%	2.276%
96	24.368	3613	3618	3629	rVB3	375158	942666	9.22%	1.312%
97	25.586	3820	3825	3838	rVB	725414	1862427	18.21%	2.592%
98	26.245	3931	3937	3953	rBV2	439612	1394958	13.64%	1.942%
99	27.621	4162	4171	4196	rVB3	1494268	6080339	59.44%	8.464%
100	28.433	4297	4309	4324	rBV3	2379077	10228525	100.00%	14.238%

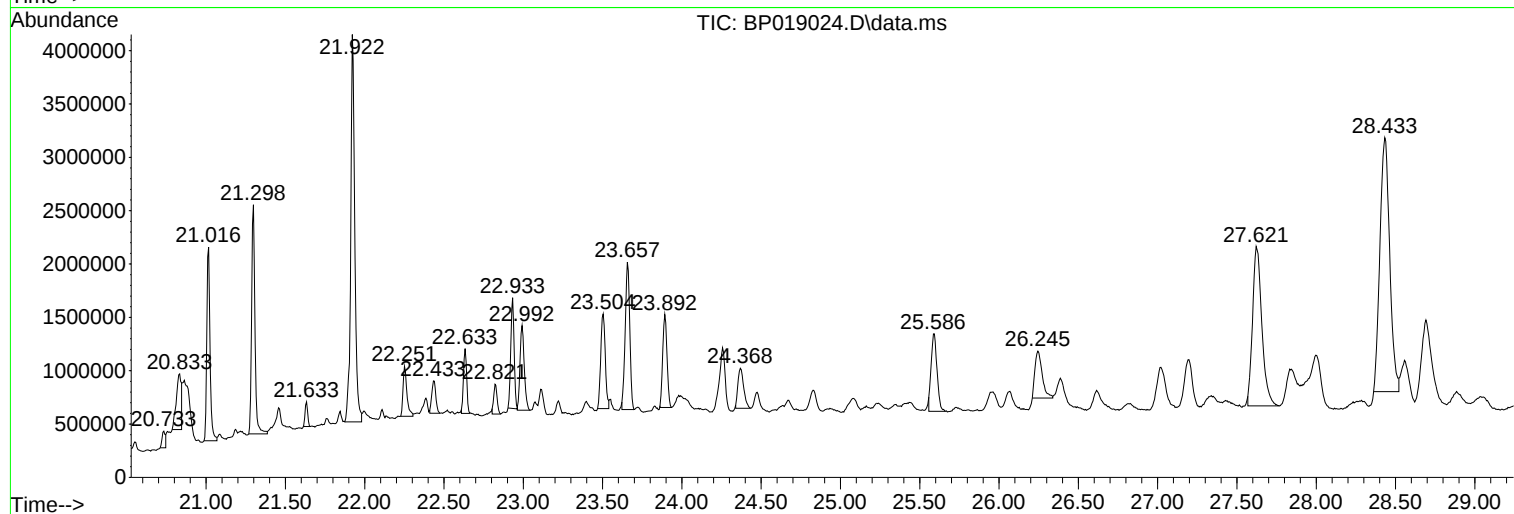
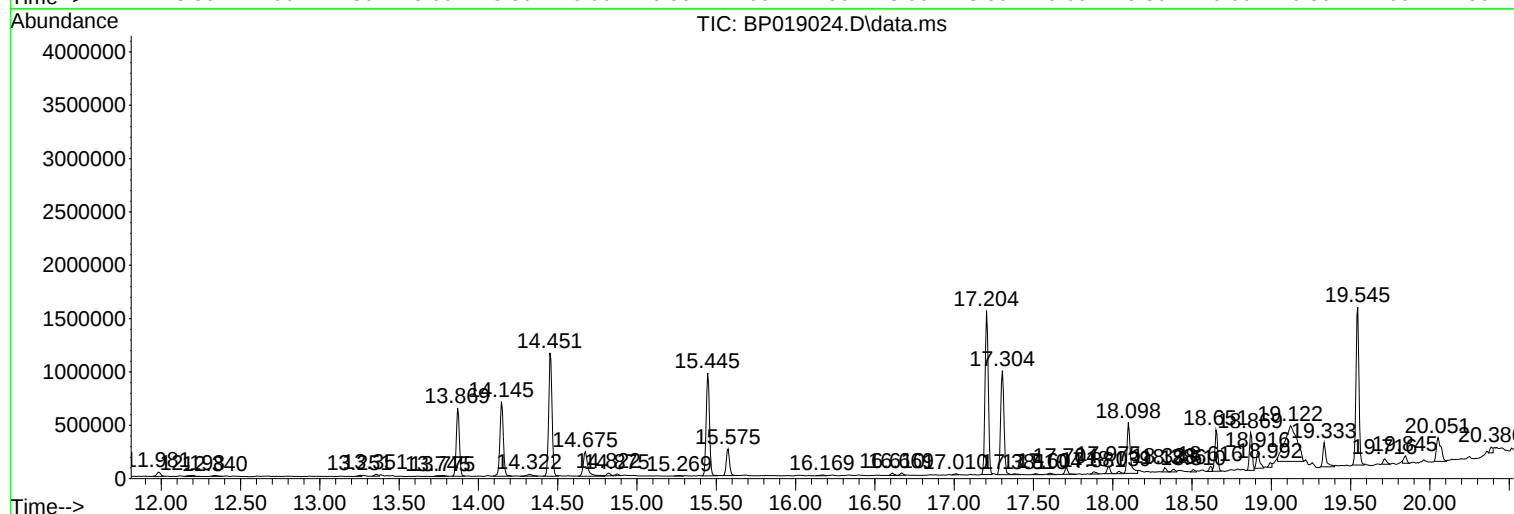
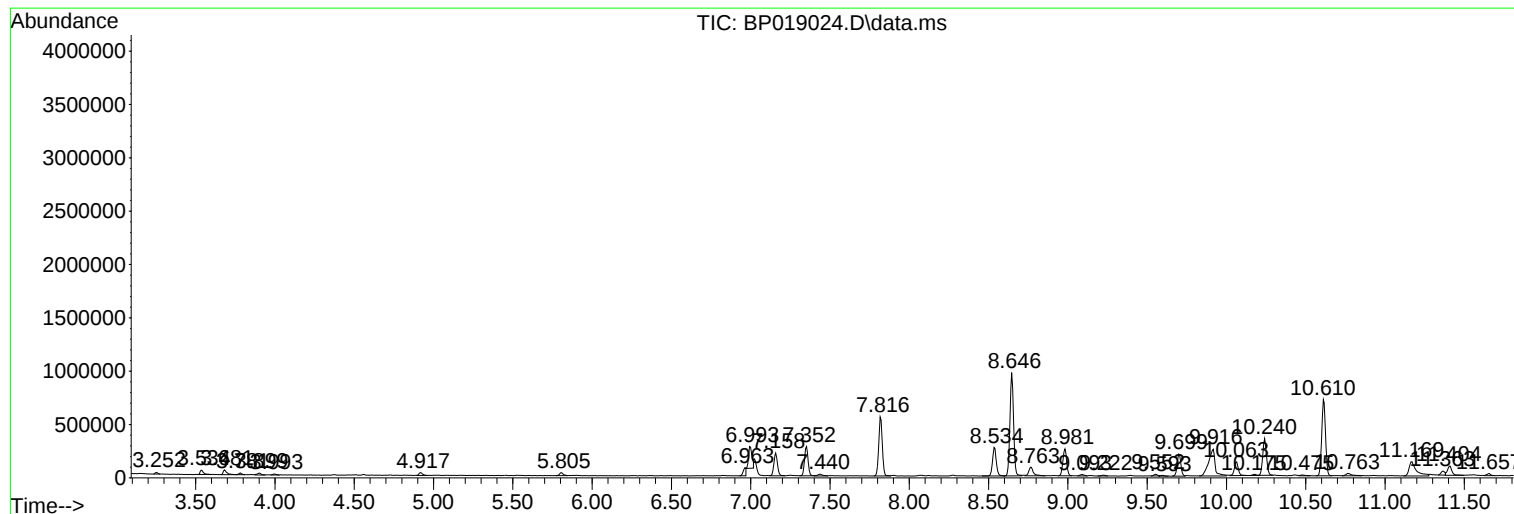
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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



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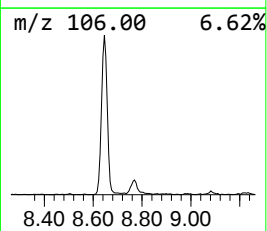
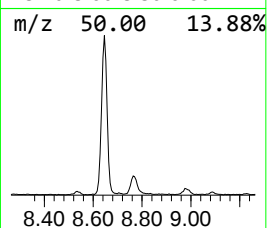
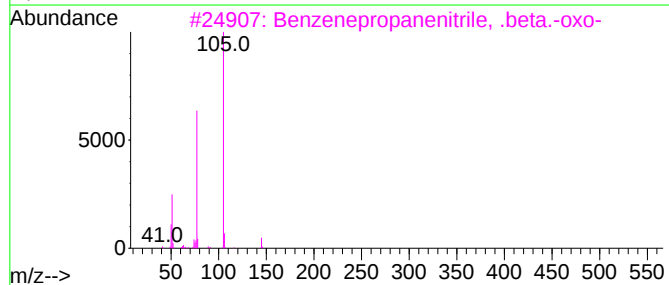
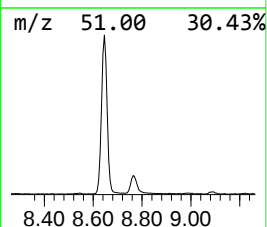
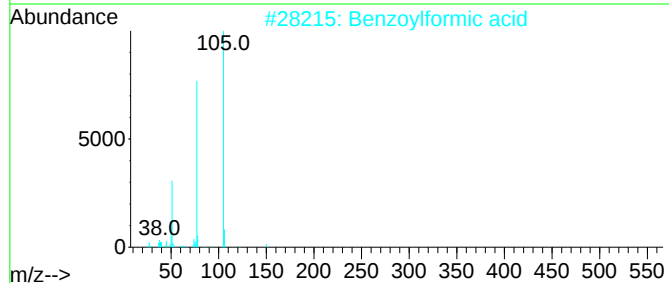
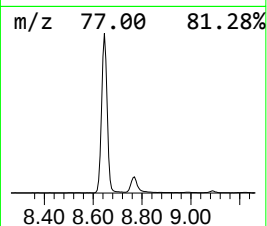
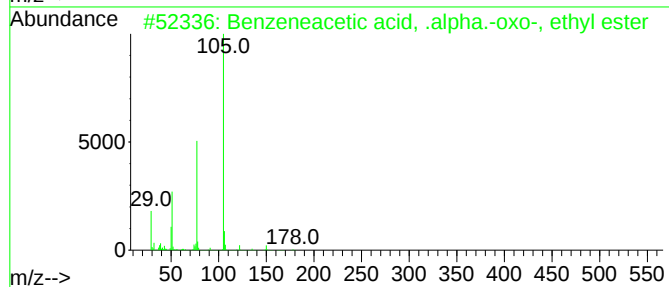
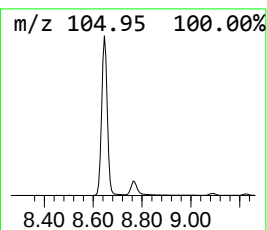
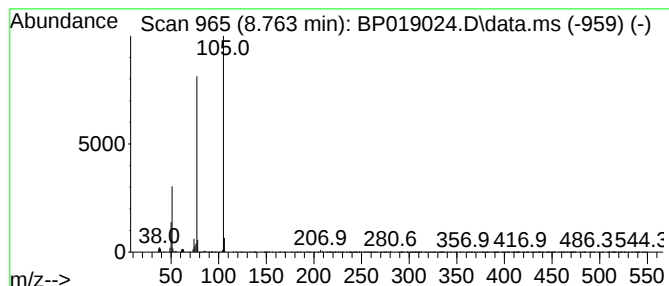
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzeneacetic acid, .alpha.... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.763	3.72 ng/ul	161325	1,4-Dichlorobenzene-d4	7.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzeneacetic acid, .alpha.-oxo-...	178	C10H10O3	001603-79-8	83
2			Benzoylformic acid	150	C8H6O3	000611-73-4	83
3			Benzenepropanenitrile, .beta.-oxo-	145	C9H7NO	000614-16-4	83
4			N-Methoxy-N-methylbenzamide	165	C9H11NO2	006919-61-5	83
5			Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	83



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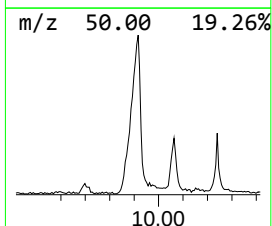
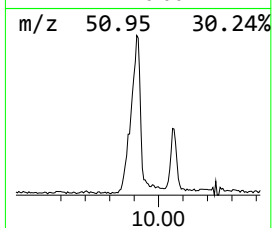
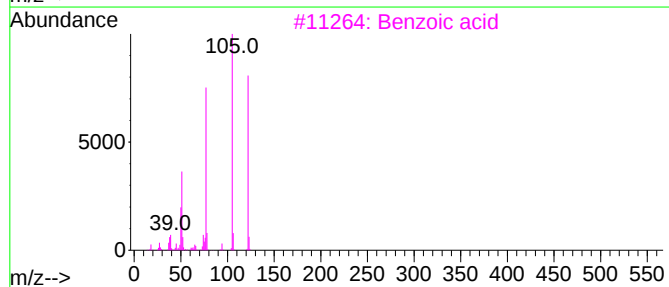
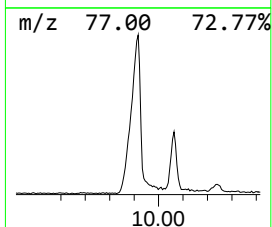
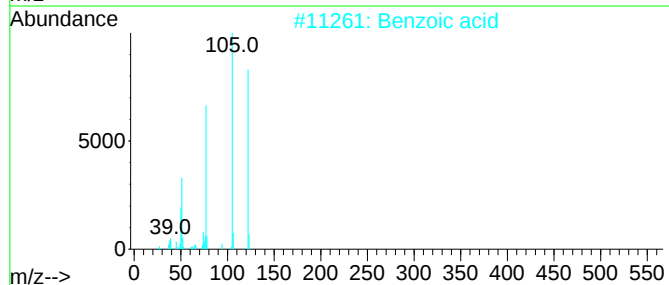
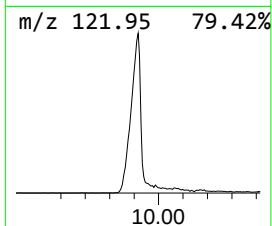
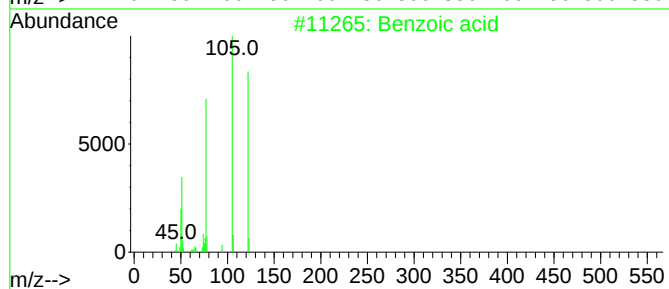
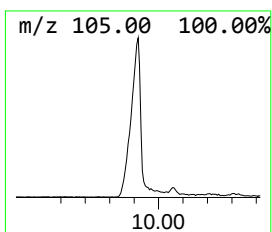
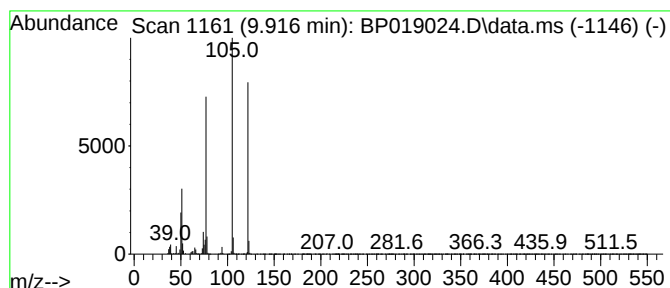
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.916	11.97 ng/ul	712552	Naphthalene-d8	10.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid			122	C7H6O2	000065-85-0	96
2	Benzoic acid			122	C7H6O2	000065-85-0	94
3	Benzoic acid			122	C7H6O2	000065-85-0	93
4	Benzoic acid			122	C7H6O2	000065-85-0	90
5	Benzoic acid, silver(1+) salt			228	C7H5AgO2	000532-31-0	90



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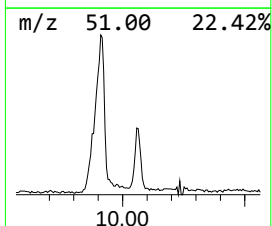
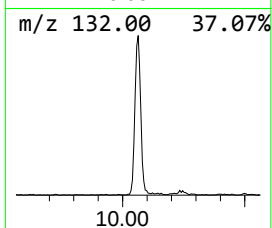
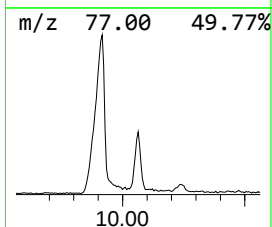
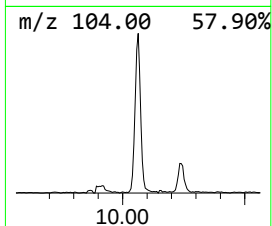
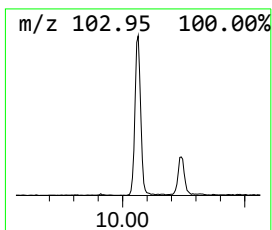
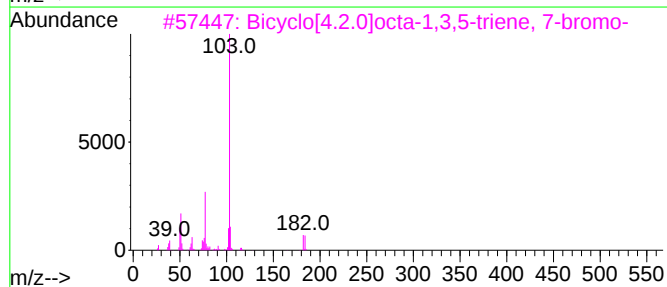
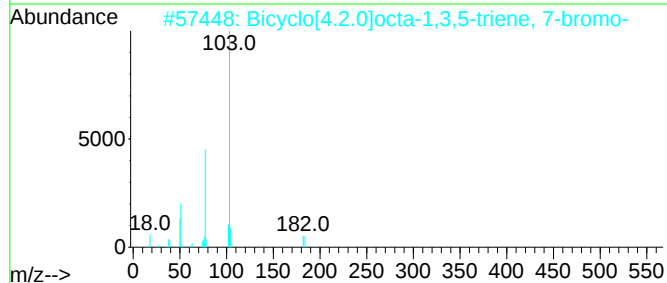
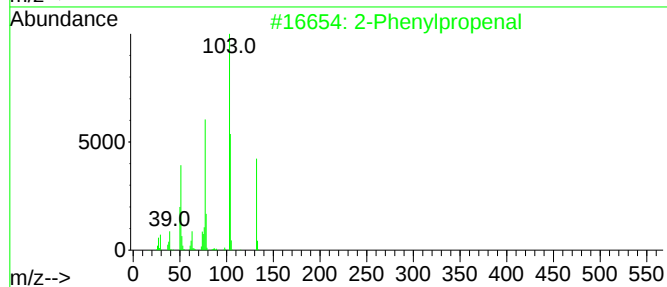
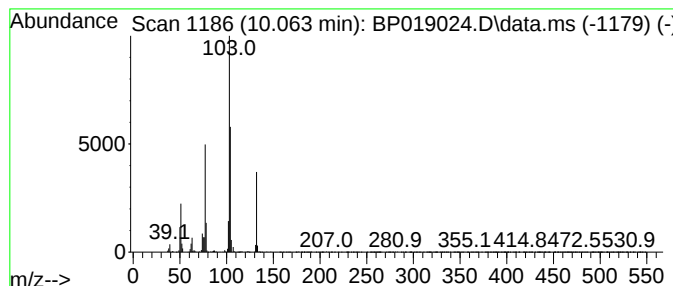
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 2-Phenylpropenal Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.063	3.70 ng/ul	220493	Naphthalene-d8	10.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Phenylpropenal	132	C9H8O	004432-63-7	70
2			Bicyclo[4.2.0]octa-1,3,5-triene,...	182	C8H7Br	021120-91-2	53
3			Bicyclo[4.2.0]octa-1,3,5-triene,...	182	C8H7Br	021120-91-2	50
4			Benzene, (1-bromoethenyl)-	182	C8H7Br	000098-81-7	50
5			Benzene, (1-iodoethenyl)-	230	C8H7I	051246-20-9	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019024.D
Acq On : 22 Dec 2023 05:48
Operator : MA/JU
Sample : 05808-19
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0B15

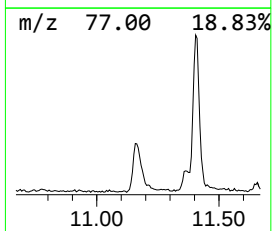
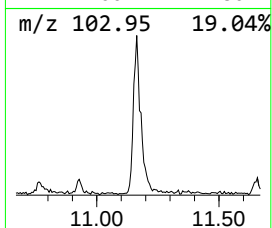
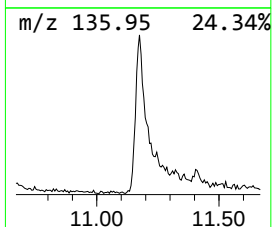
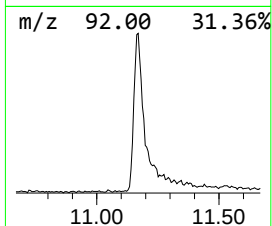
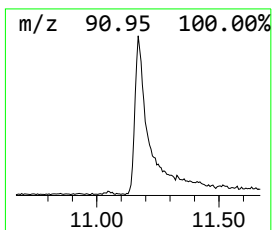
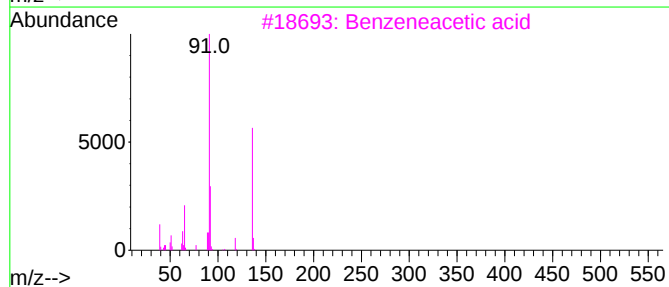
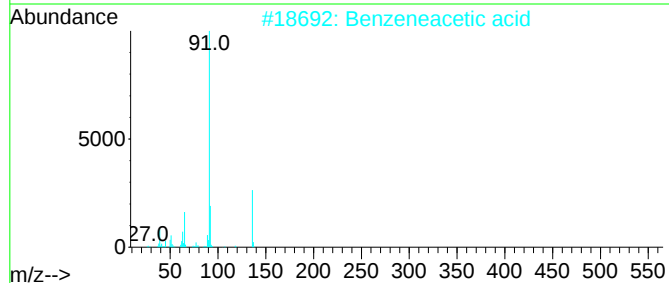
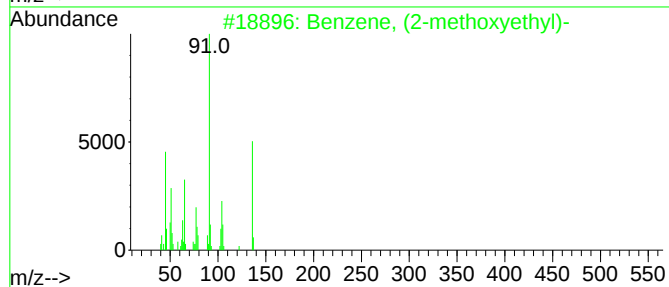
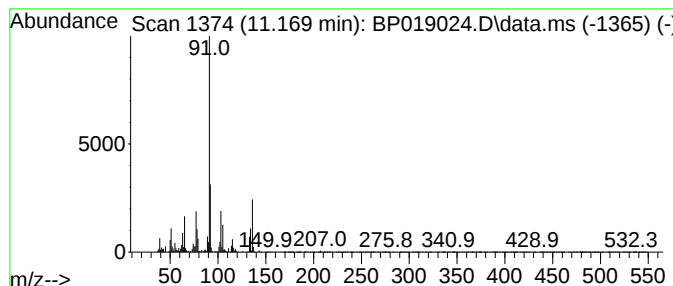
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-01 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.169	7.11 ng/ul	423241	Naphthalene-d8	10.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methoxyethyl)-	136	C9H12O	003558-60-9	49
2			Benzeneacetic acid	136	C8H8O2	000103-82-2	46
3			Benzeneacetic acid	136	C8H8O2	000103-82-2	43
4			Benzeneacetic acid	136	C8H8O2	000103-82-2	41
5			Benzeneacetic acid	136	C8H8O2	000103-82-2	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019024.D
Acq On : 22 Dec 2023 05:48
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Sample : 05808-19
Misc :
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Instrument :
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ClientSampleId :
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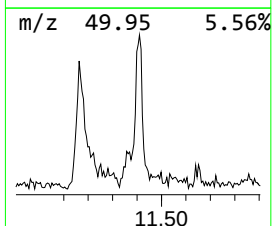
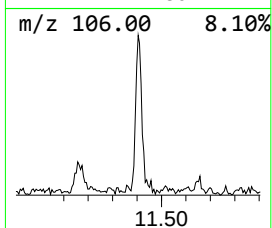
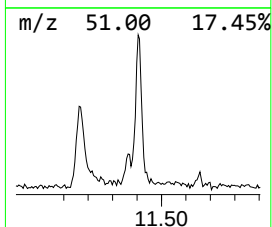
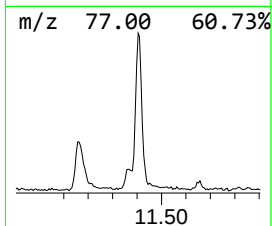
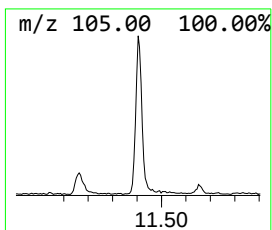
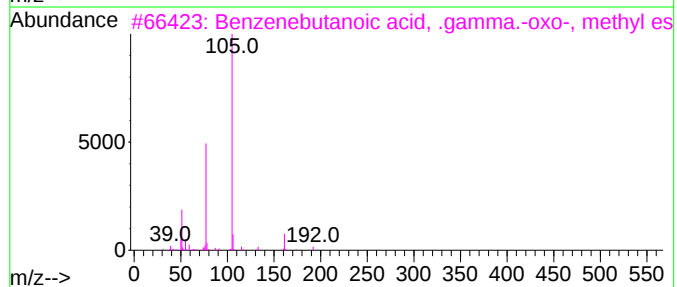
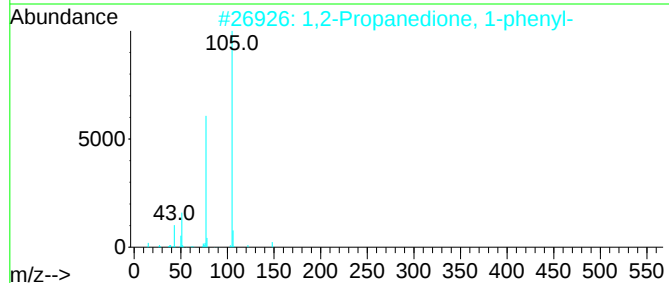
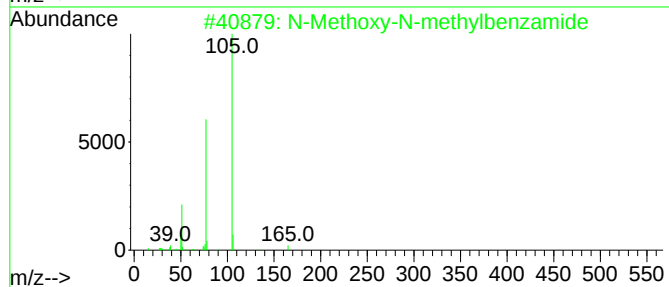
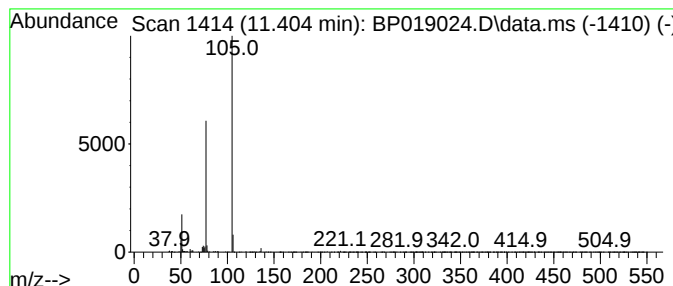
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 N-Methoxy-N-methylbenzamide Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.404	2.77 ng/ul	164649	Naphthalene-d8	10.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			N-Methoxy-N-methylbenzamide	165	C9H11NO2	006919-61-5	78
2			1,2-Propanedione, 1-phenyl-	148	C9H8O2	000579-07-7	78
3			Benzenebutanoic acid, .gamma.-ox...	192	C11H12O3	025333-24-8	78
4			1,2-Propanedione, 1-phenyl-	148	C9H8O2	000579-07-7	78
5			Benzeneacetic acid, .alpha.-oxo-...	164	C9H8O3	015206-55-0	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019024.D
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Operator : MA/JU
Sample : 05808-19
Misc :
ALS Vial : 24 Sample Multiplier: 1

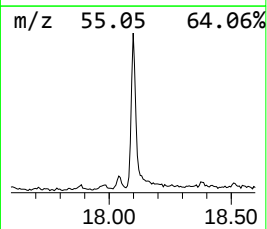
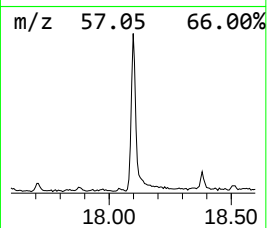
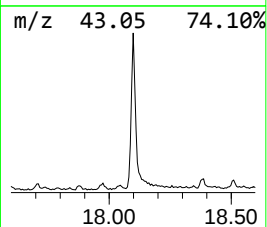
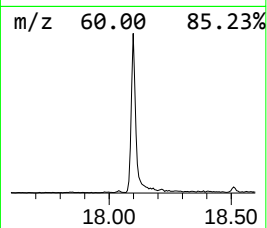
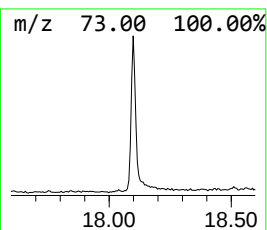
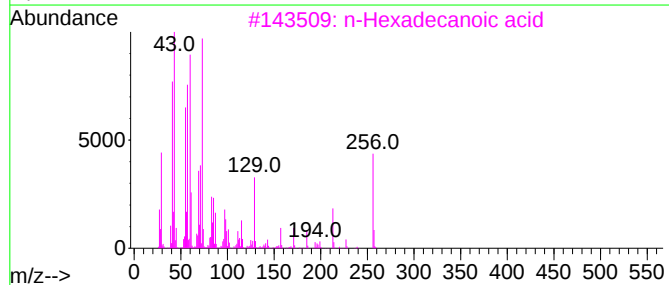
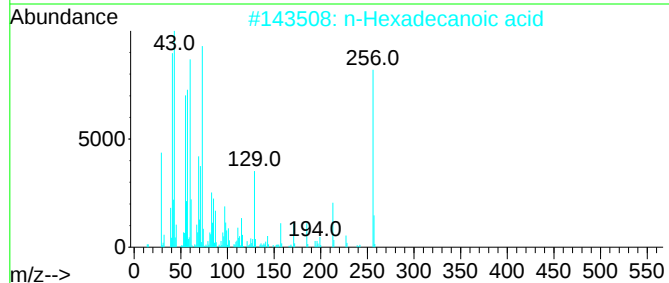
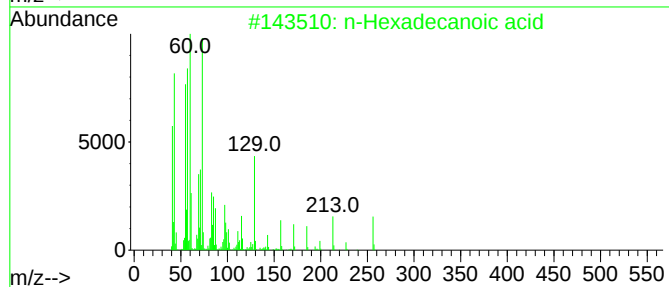
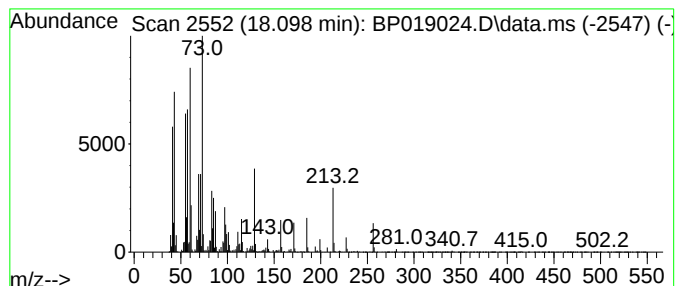
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.098	6.35 ng/ul	679229	Phenanthrene-d10	17.204	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4	Tridecanoic acid	214	C13H26O2	000638-53-9	97
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019024.D
Acq On : 22 Dec 2023 05:48
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Sample : 05808-19
Misc :
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Instrument :
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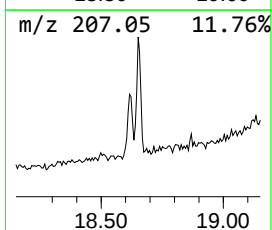
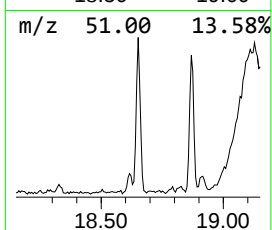
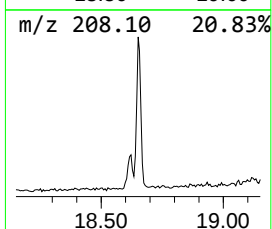
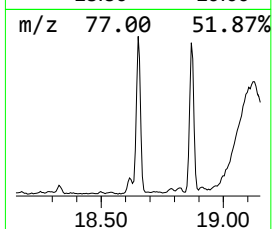
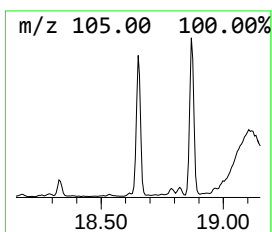
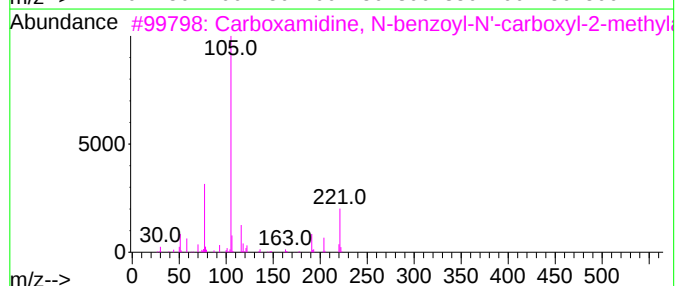
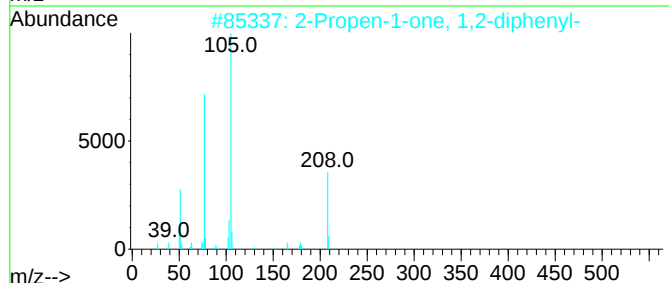
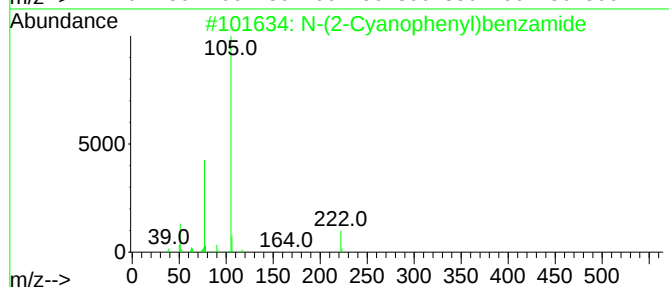
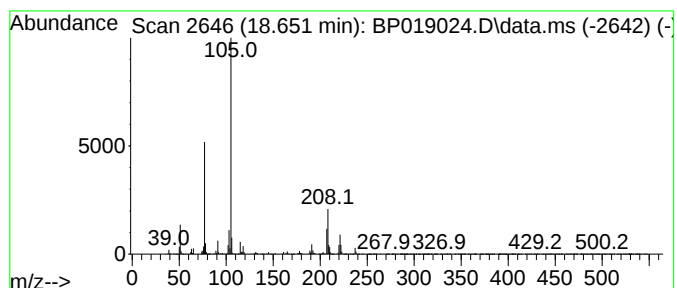
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 N-(2-Cyanophenyl)benzamide Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.651	4.43 ng/ul	473494	Phenanthrene-d10	17.204

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			N-(2-Cyanophenyl)benzamide	222	C14H10N2O	040288-69-5	50
2			2-Propen-1-one, 1,2-diphenyl-	208	C15H12O	004452-11-3	49
3			Carboximidine, N-benzoyl-N'-carb...	221	C10H11N3O3	1000160-14-9	47
4			Ethanone, 2-(2-ethenylphenyl)-1-...	222	C16H14O	066657-90-7	45
5			Benzamide, N-(3-ethynylphenyl)-	221	C15H11NO	208943-24-2	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019024.D
Acq On : 22 Dec 2023 05:48
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Sample : 05808-19
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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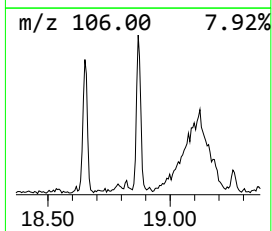
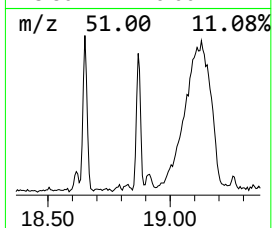
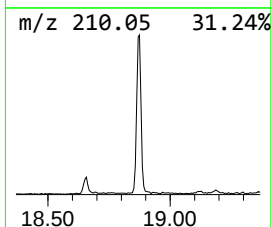
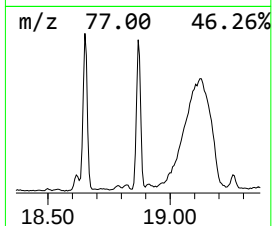
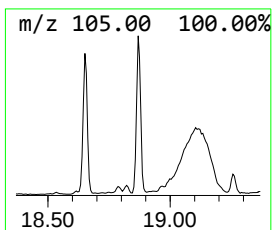
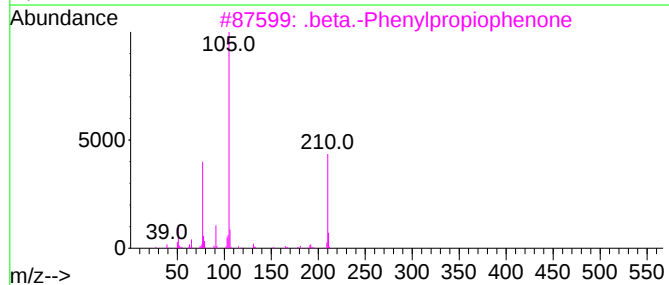
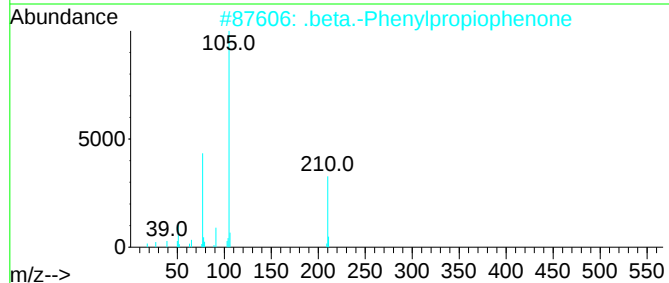
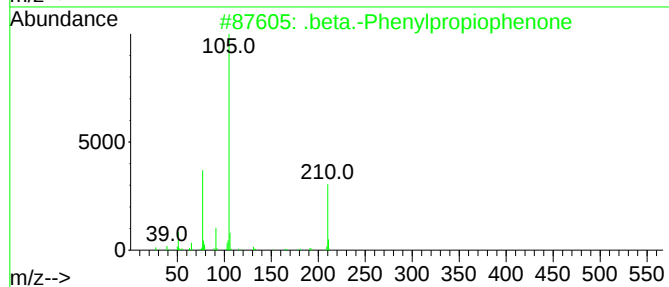
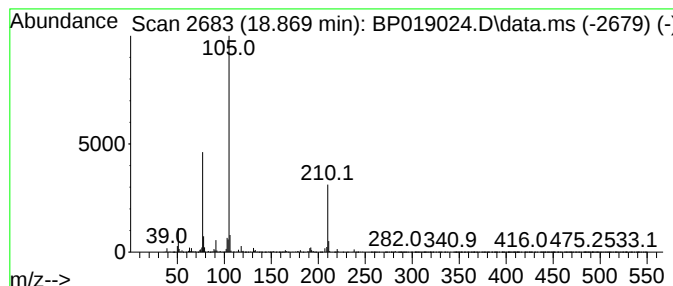
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 .beta.-Phenylpropiophenone Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.869	4.03 ng/ul	431238	Phenanthrene-d10	17.204

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.	.beta.-Phenylpropiophenone	210	C15H14O	001083-30-3	93	
2	.	.beta.-Phenylpropiophenone	210	C15H14O	001083-30-3	93	
3	.	.beta.-Phenylpropiophenone	210	C15H14O	001083-30-3	89	
4	.	.beta.-Phenylpropiophenone	210	C15H14O	001083-30-3	87	
5	N-Benzoyl-3-phenylisoserine, eth...	341	C20H23NO4	1010453-73-5	72		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019024.D
Acq On : 22 Dec 2023 05:48
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Misc :
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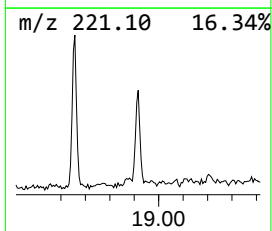
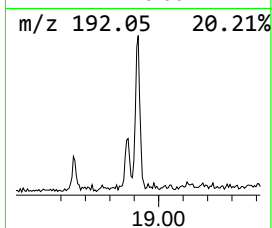
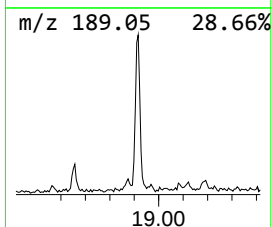
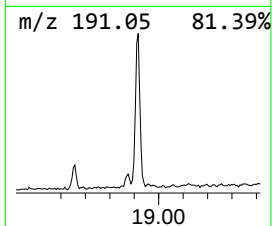
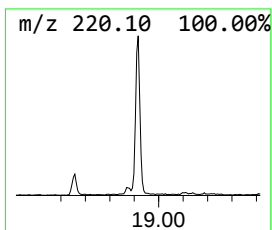
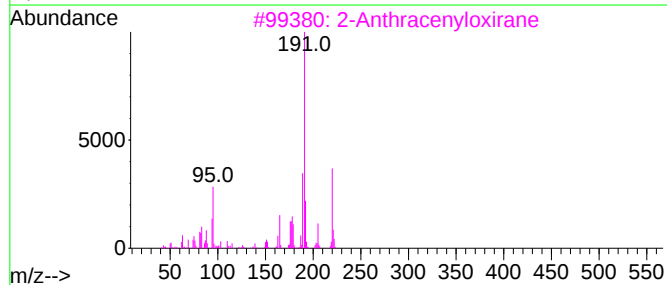
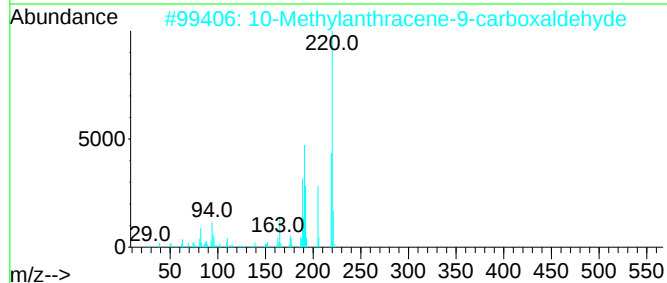
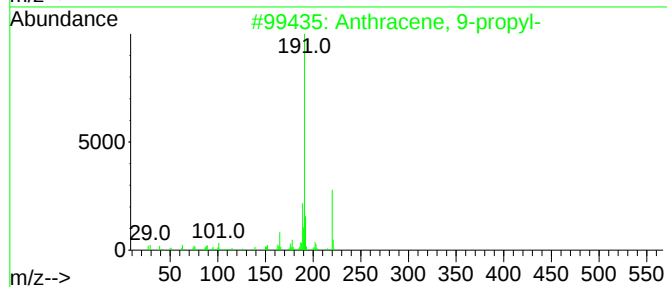
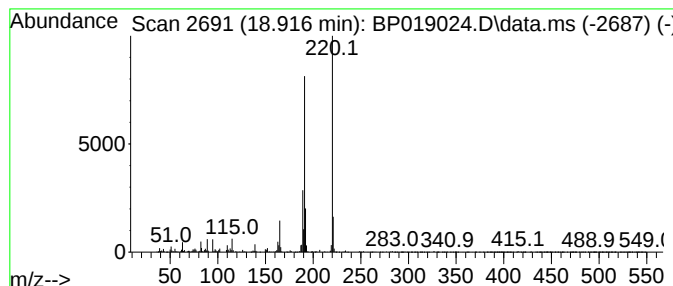
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Anthracene, 9-propyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.916	2.06 ng/ul	219938	Phenanthrene-d10	17.204

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 9-propyl-	220	C17H16	001498-77-7	87
2			10-Methylanthracene-9-carboxalde...	220	C16H12O	007072-00-6	86
3			2-Anthracenyloxirane	220	C16H12O	052643-86-4	72
4			1H-Pyrazole, 3,5-diphenyl-	220	C15H12N2	001145-01-3	58
5			1-(9-Phenanthrenyl)-2-propanone	234	C17H14O	1000326-08-9	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019024.D
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Operator : MA/JU
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Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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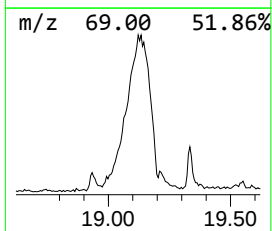
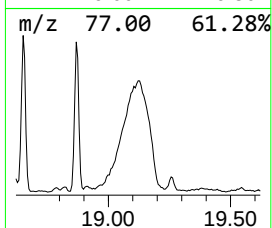
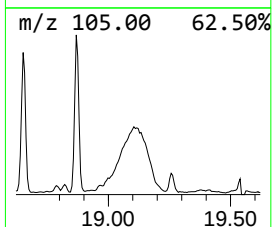
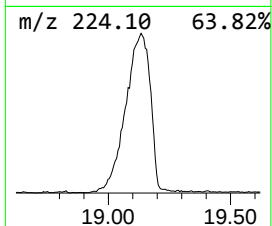
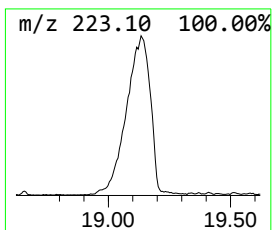
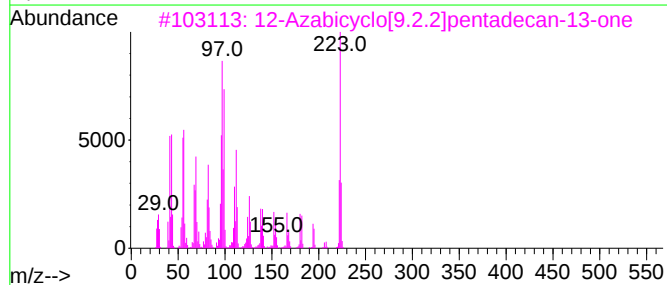
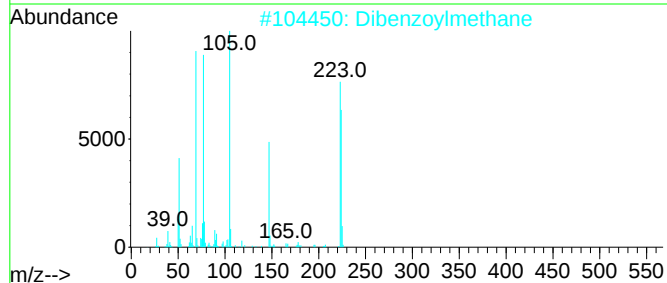
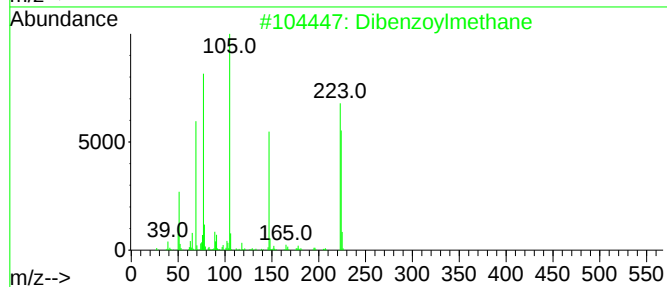
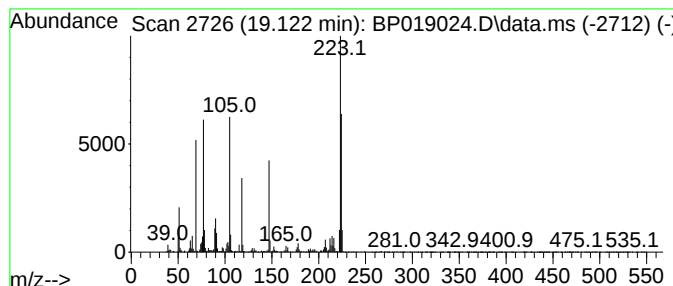
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Dibenzoylmethane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.122	16.73 ng/ul	1790210	Phenanthrene-d10	17.204

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzoylmethane	224	C15H12O2	000120-46-7	97
2			Dibenzoylmethane	224	C15H12O2	000120-46-7	96
3			12-Azabicyclo[9.2.2]pentadecan-1...	223	C14H25NO	1000191-26-7	94
4			Dibenzoylmethane	224	C15H12O2	000120-46-7	91
5			Dibenzoylmethane	224	C15H12O2	000120-46-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019024.D
Acq On : 22 Dec 2023 05:48
Operator : MA/JU
Sample : 05808-19
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0B15

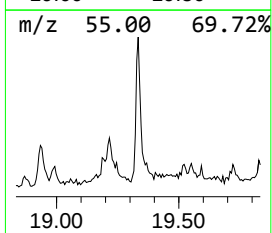
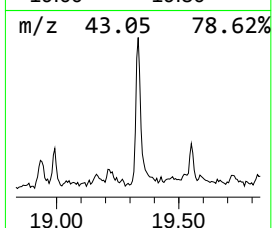
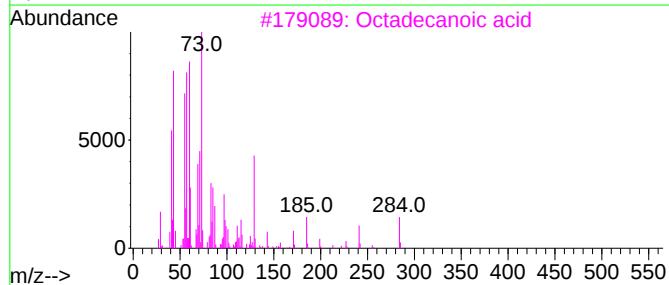
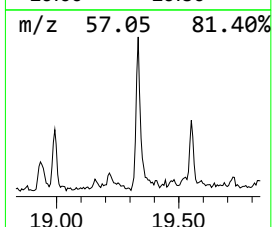
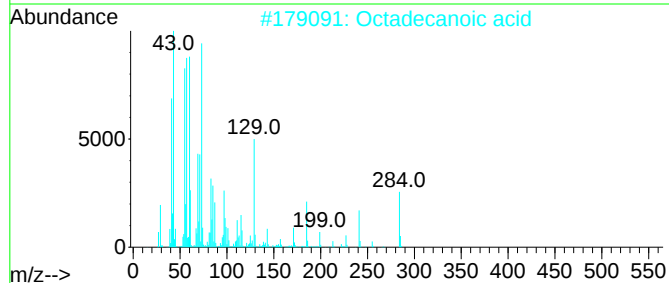
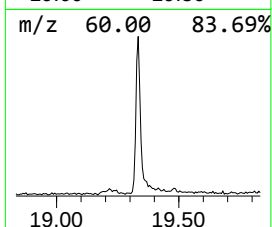
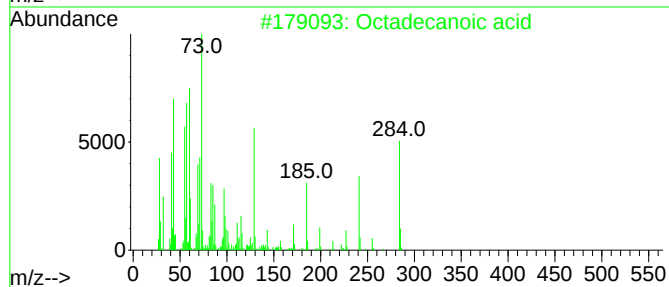
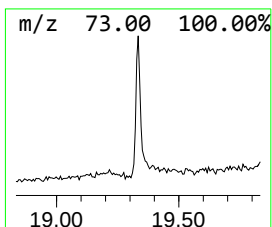
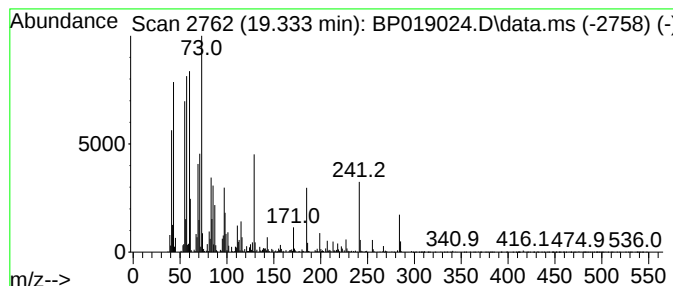
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Octadecanoic acid Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.333	2.05 ng/ul	324457	Chrysene-d12	21.298

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	98
3			Octadecanoic acid	284	C18H36O2	000057-11-4	98
4			Octadecanoic acid	284	C18H36O2	000057-11-4	95
5			Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019024.D
Acq On : 22 Dec 2023 05:48
Operator : MA/JU
Sample : 05808-19
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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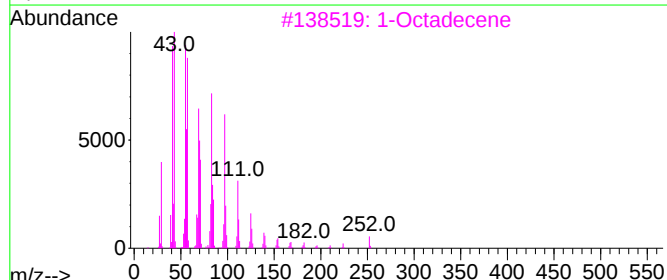
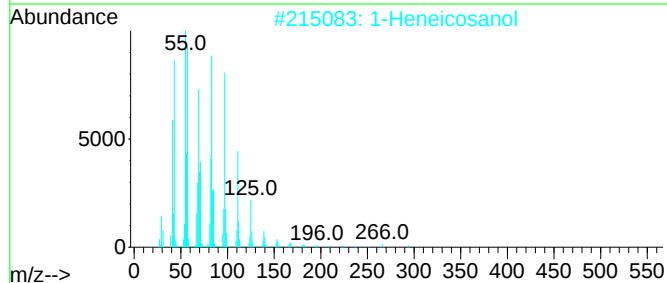
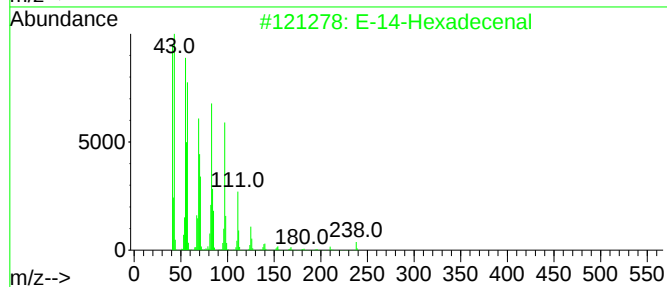
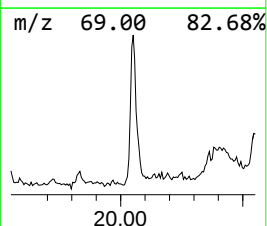
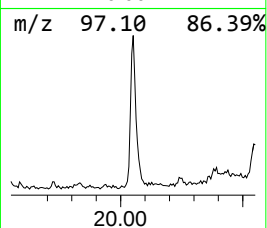
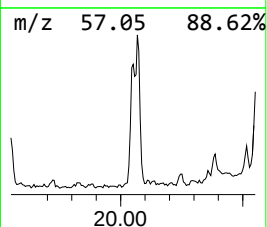
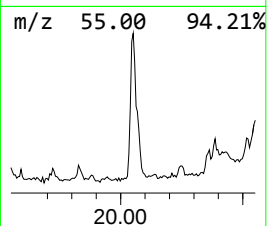
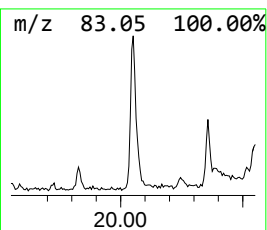
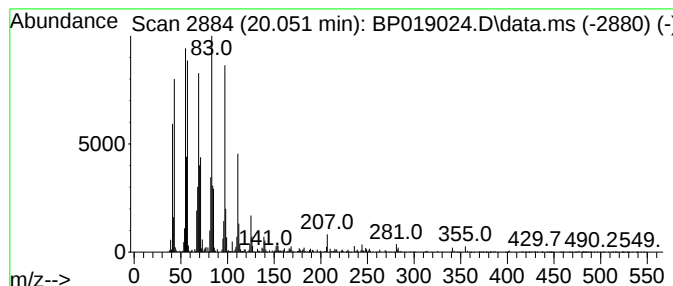
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 E-14-Hexadecenal Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.051	2.78 ng/ul	440838	Chrysene-d12	21.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	E-14-Hexadecenal			238	C16H30O	330207-53-9	98
2	1-Heneicosanol			312	C21H44O	015594-90-8	95
3	1-Octadecene			252	C18H36	000112-88-9	94
4	n-Nonadecanol-1			284	C19H40O	001454-84-8	94
5	n-Nonadecanol-1			284	C19H40O	001454-84-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019024.D
Acq On : 22 Dec 2023 05:48
Operator : MA/JU
Sample : 05808-19
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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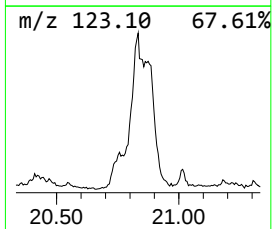
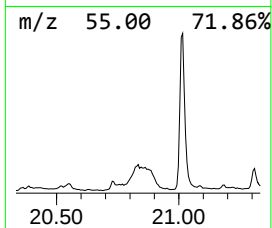
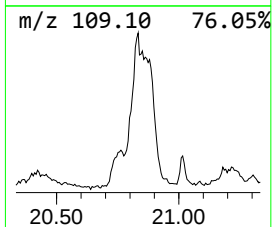
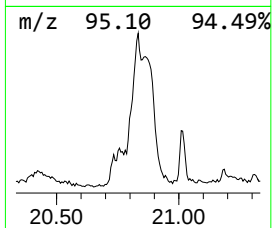
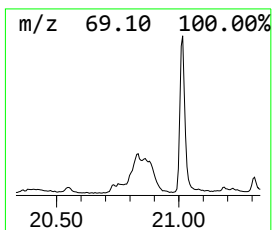
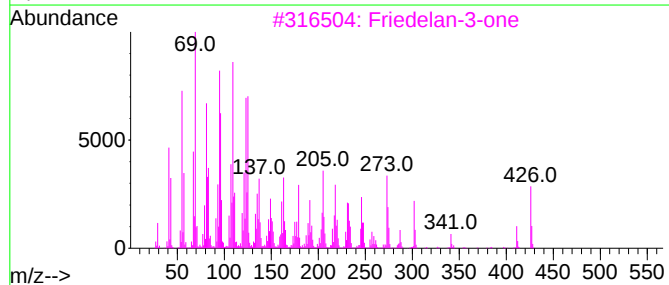
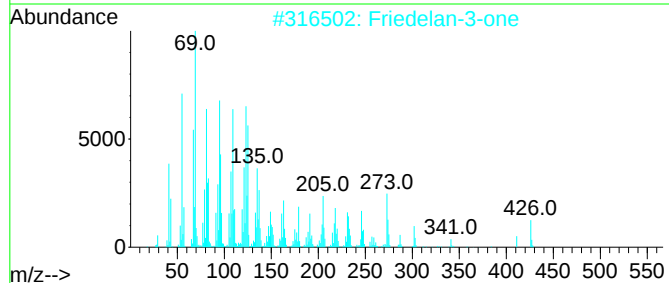
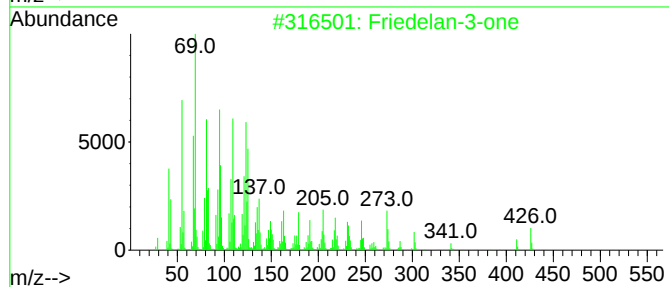
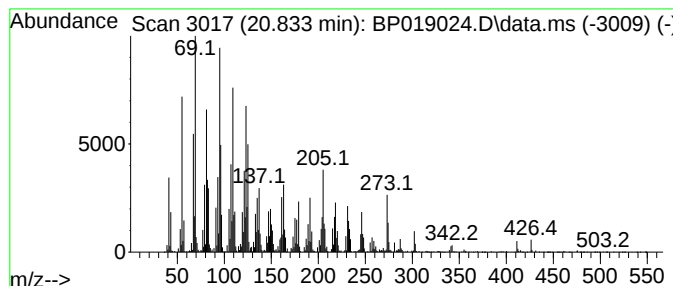
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 5-Bromo-4-oxo-4,5,6,7-tetra... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.833	7.28 ng/ul	1153050	Chrysene-d12	21.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Friedelan-3-one	426	C30H50O	000559-74-0	95
2			Friedelan-3-one	426	C30H50O	000559-74-0	91
3			Friedelan-3-one	426	C30H50O	000559-74-0	78
4			5-Bromo-4-oxo-4,5,6,7-tetrahydro...	216	C6H5BrN2O2	300574-36-1	60
5			(E)-3-Methyl-5-((1R,4aR,8aR)-5,5...	290	C20H34O	021738-29-4	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019024.D
Acq On : 22 Dec 2023 05:48
Operator : MA/JU
Sample : 05808-19
Misc :
ALS Vial : 24 Sample Multiplier: 1

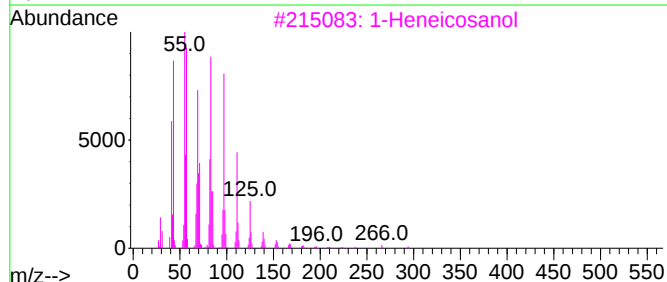
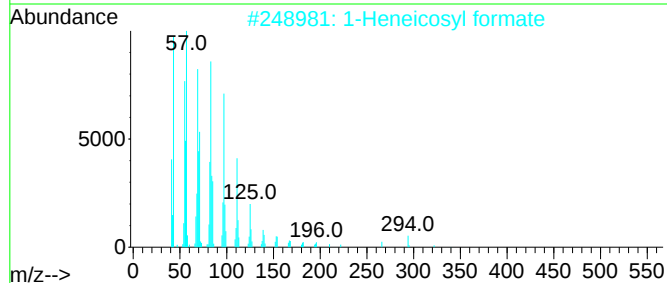
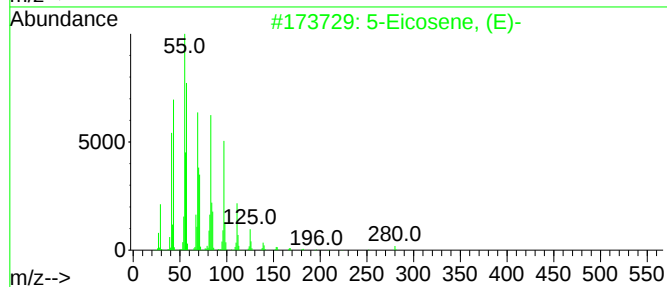
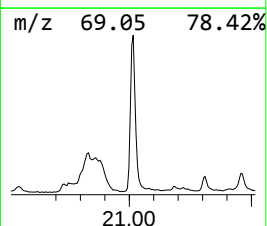
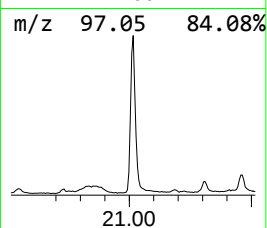
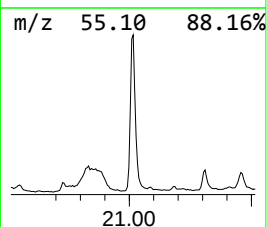
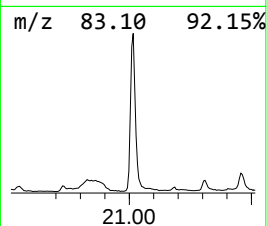
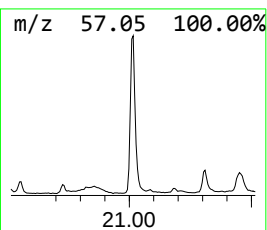
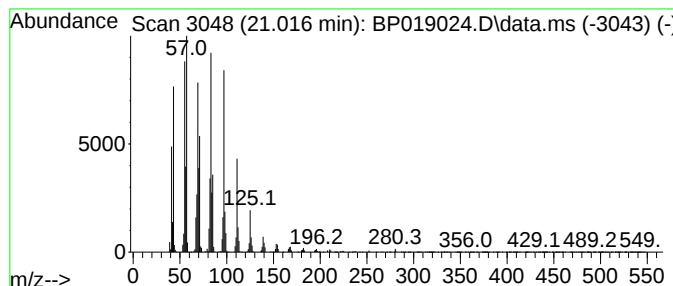
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.016	16.21 ng/ul	2566950	Chrysene-d12		21.298	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	5-Eicosene, (E)-		280	C20H40	074685-30-6	98
2	1-Heneicosyl formate		340	C22H44O2	077899-03-7	95
3	1-Heneicosanol		312	C21H44O	015594-90-8	95
4	Behenic alcohol		326	C22H46O	000661-19-8	94
5	Heptadecyl heptafluorobutyrate		452	C21H35F7O2	959085-66-6	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019024.D
Acq On : 22 Dec 2023 05:48
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Sample : 05808-19
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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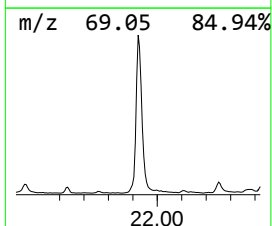
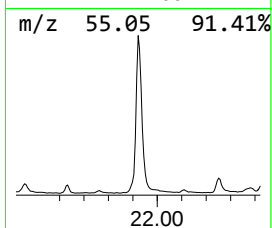
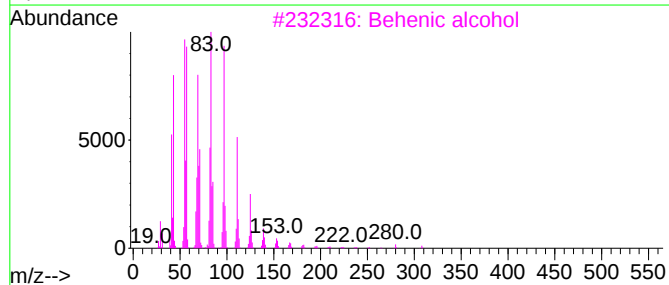
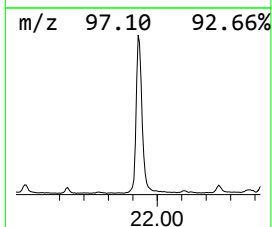
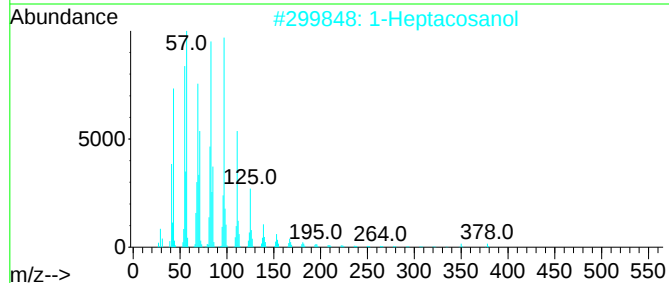
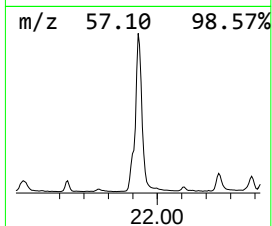
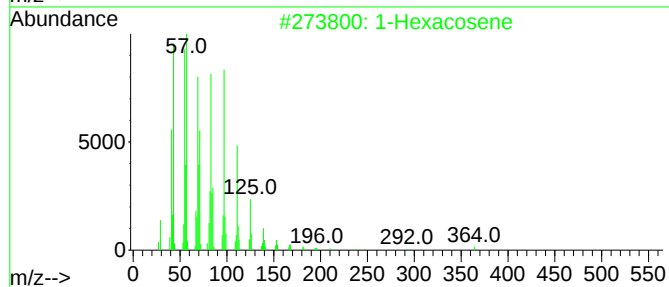
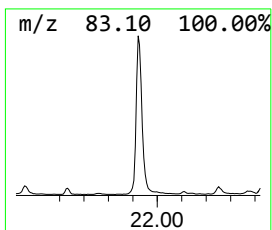
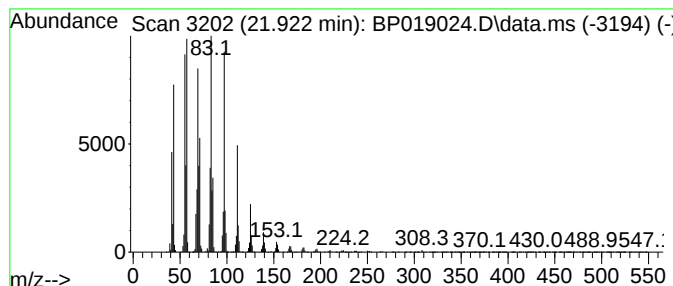
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.922	41.18 ng/ul	6522120	Chrysene-d12	21.298

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexacosene		364	C26H52	018835-33-1	95
2	1-Heptacosanol		396	C27H56O	002004-39-9	95
3	Behenic alcohol		326	C22H46O	000661-19-8	95
4	1-Heneicosanol		312	C21H44O	015594-90-8	94
5	n-Tetracosanol-1		354	C24H50O	000506-51-4	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019024.D
Acq On : 22 Dec 2023 05:48
Operator : MA/JU
Sample : 05808-19
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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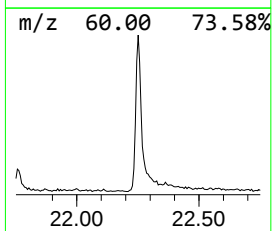
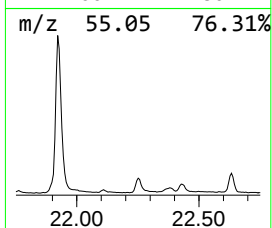
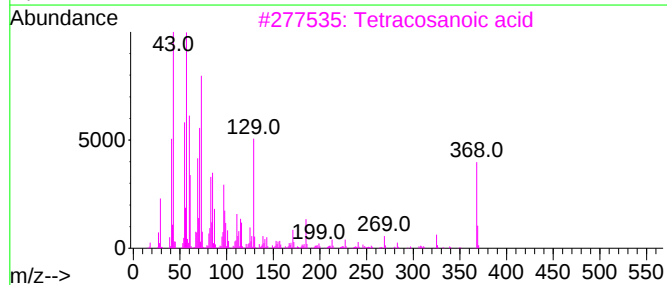
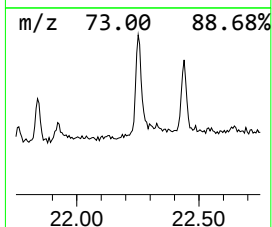
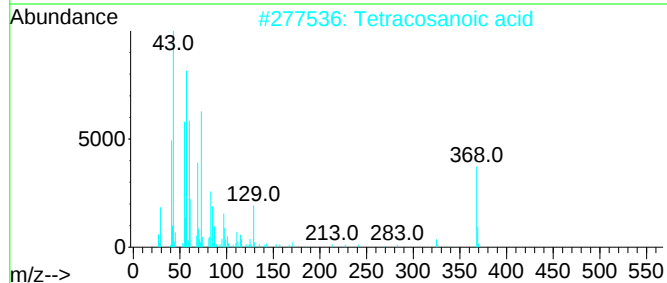
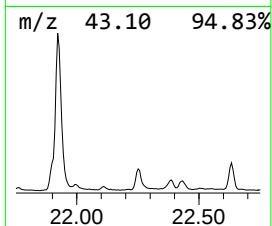
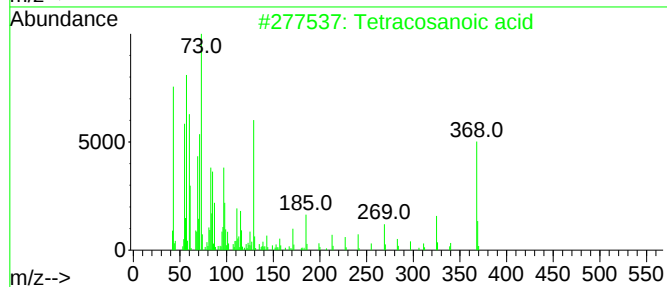
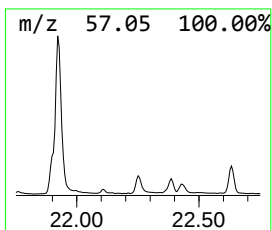
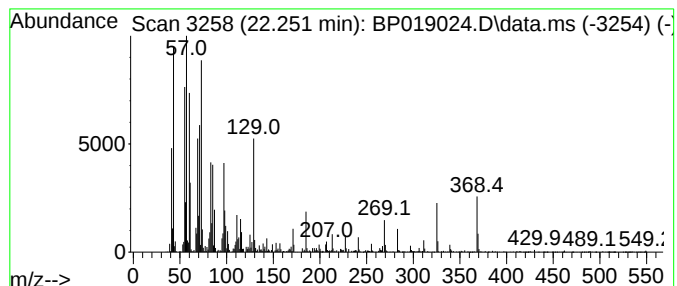
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 Tetracosanoic acid Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.251	4.97 ng/ul	787338	Chrysene-d12	21.298

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosanoic acid	368	C24H48O2	000557-59-5	99
2		Tetracosanoic acid	368	C24H48O2	000557-59-5	99
3		Tetracosanoic acid	368	C24H48O2	000557-59-5	92
4		n-Decanoic acid	172	C10H20O2	000334-48-5	89
5		Heneicosanoic acid	326	C21H42O2	002363-71-5	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019024.D
Acq On : 22 Dec 2023 05:48
Operator : MA/JU
Sample : 05808-19
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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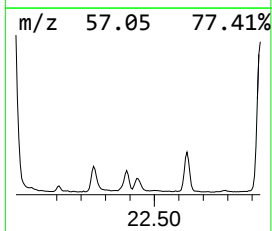
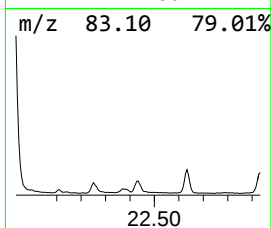
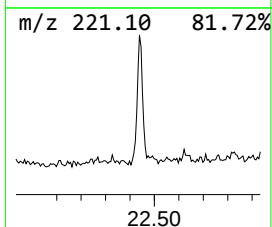
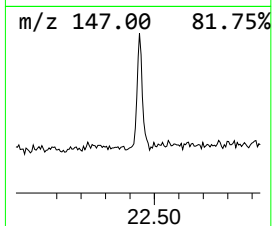
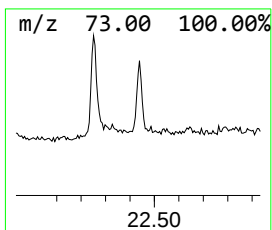
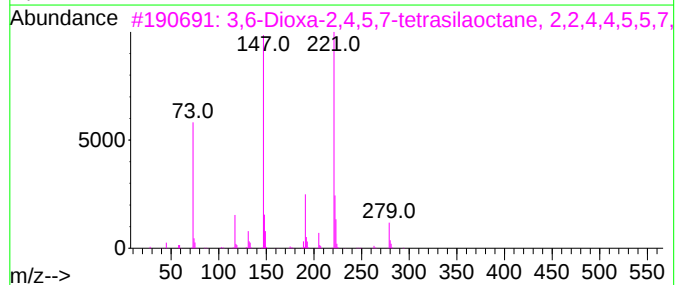
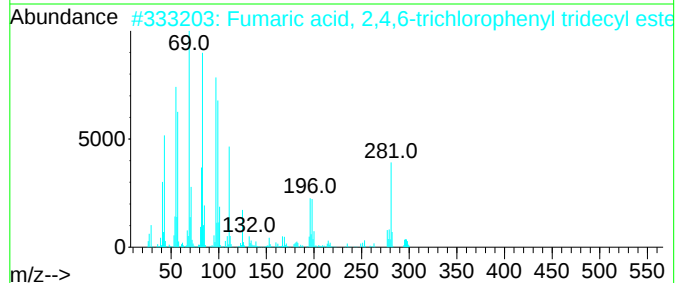
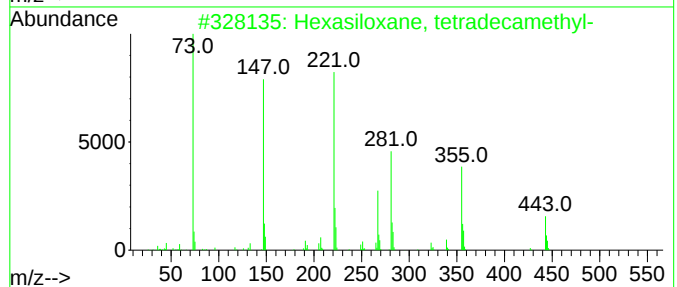
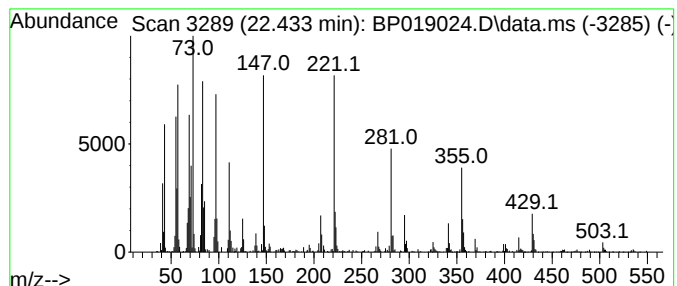
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 unknown-02 Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.433	3.39 ng/ul	536274	Chrysene-d12	21.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexasiloxane, tetradecamethyl-	458	C14H42O5Si6	000107-52-8	46
2			Fumaric acid, 2,4,6-trichlorophe...	476	C23H31Cl3O4	1000348-27-7	27
3			3,6-Dioxa-2,4,5,7-tetrasilaoctan...	294	C10H30O2Si4	004342-25-0	25
4			Silane, [(11-iodoundecyl)oxy]tri...	370	C14H31IO5Si	026305-84-0	16
5			1,1,1,5,7,7,7-Heptamethyl-3,3-bi...	444	C13H40O5Si6	038147-00-1	16



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019024.D
Acq On : 22 Dec 2023 05:48
Operator : MA/JU
Sample : 05808-19
Misc :
ALS Vial : 24 Sample Multiplier: 1

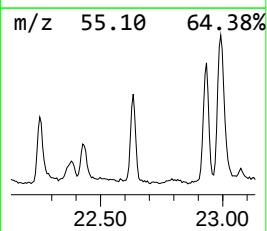
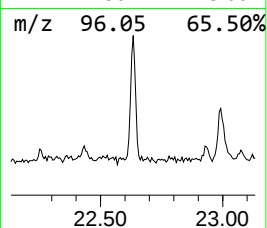
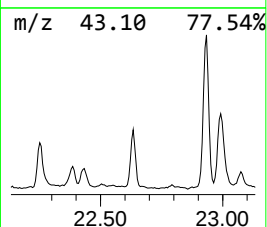
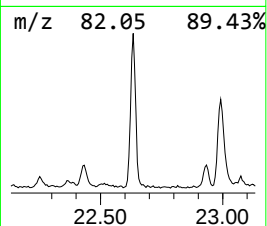
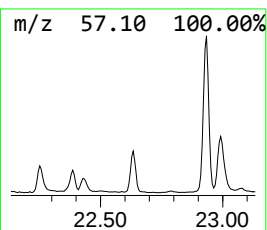
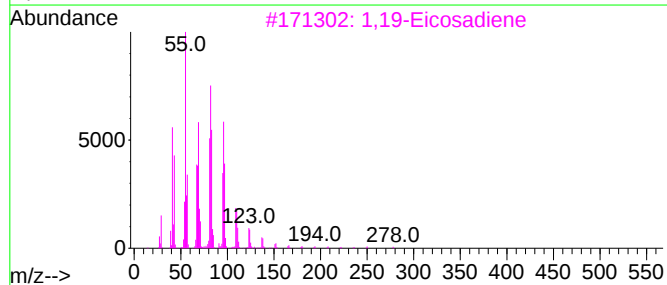
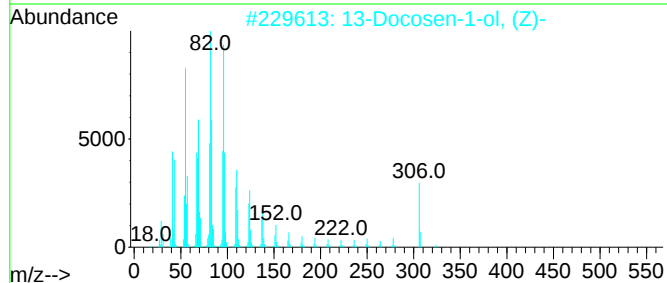
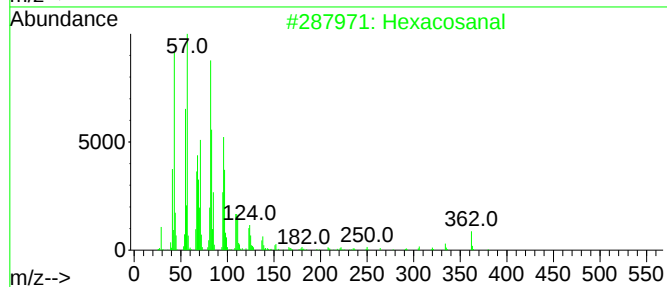
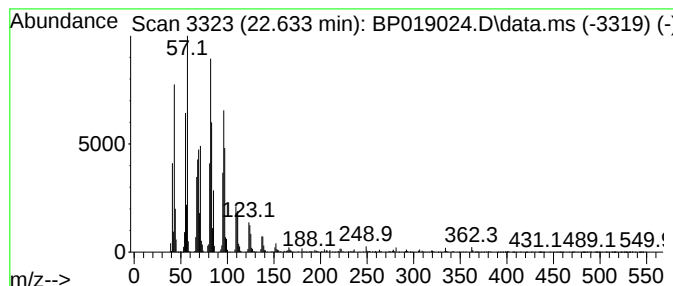
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 18 Hexacosanal Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
22.633	6.13 ng/ul	842991	Perylene-d12	23.657	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosanal	380	C26H52O	026627-85-0	97
2	13-Docosen-1-ol, (Z)-	324	C22H44O	000629-98-1	96
3	1,19-Eicosadiene	278	C20H38	014811-95-1	94
4	Oxirane, hexadecyl-	268	C18H36O	007390-81-0	94
5	Oxirane, heptadecyl-	282	C19H38O	067860-04-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019024.D
Acq On : 22 Dec 2023 05:48
Operator : MA/JU
Sample : 05808-19
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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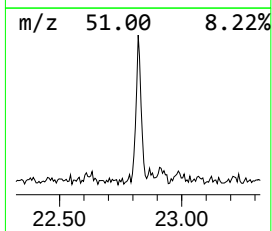
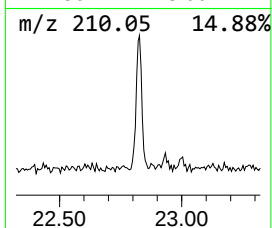
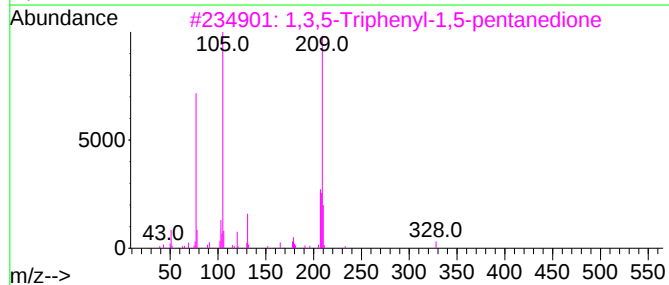
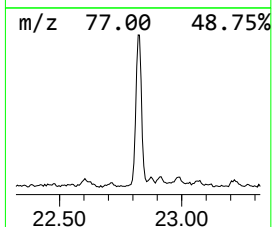
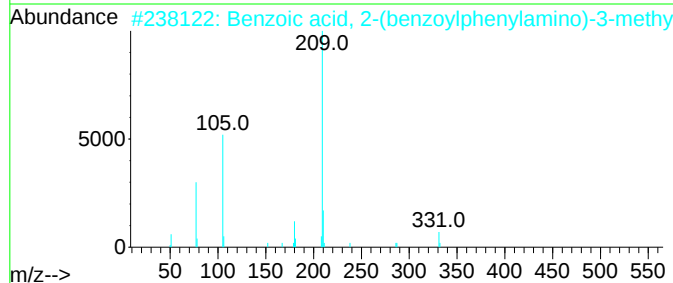
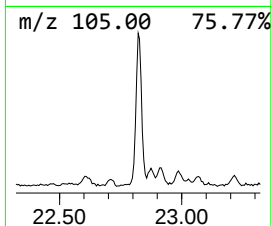
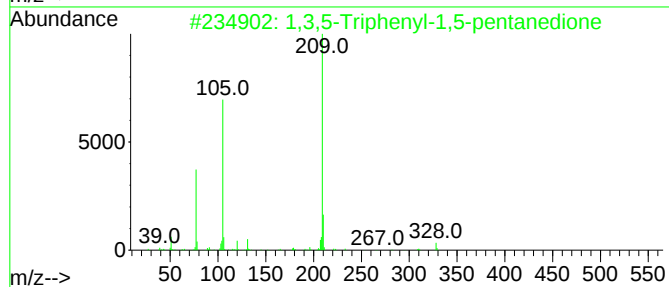
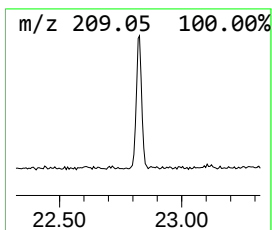
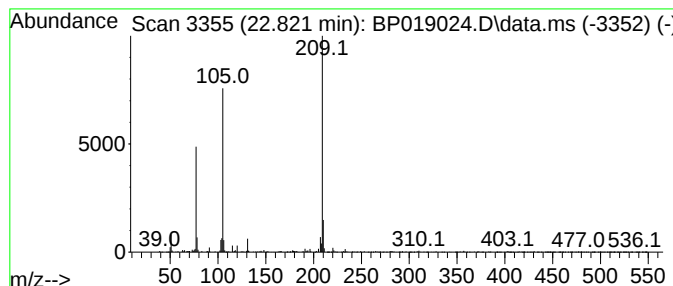
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 1,3,5-Triphenyl-1,5-pentane... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.822	2.99 ng/ul	410571	Perylene-d12	23.657

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3,5-Triphenyl-1,5-pentanedione	328	C23H20O2	006263-84-9	78	
2	Benzoic acid, 2-(benzoylphenylam...	331	C21H17NO3	050405-25-9	78	
3	1,3,5-Triphenyl-1,5-pentanedione	328	C23H20O2	006263-84-9	64	
4	1-[(4-benzoylbenzoyl)oxy] pyrrol...	323	C18H13NO5	1000334-29-5	50	
5	Benzamide, 4-benzoyl-N-(2-triflu...	369	C21H14F3NO2	345988-01-4	50	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019024.D
Acq On : 22 Dec 2023 05:48
Operator : MA/JU
Sample : 05808-19
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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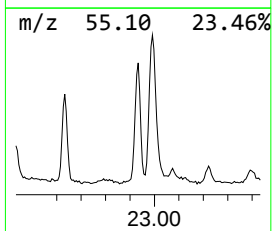
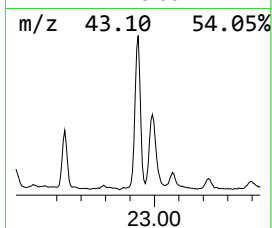
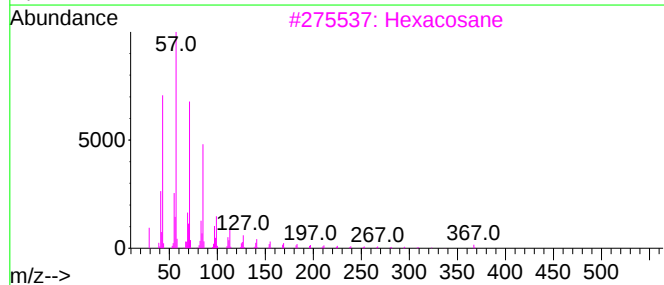
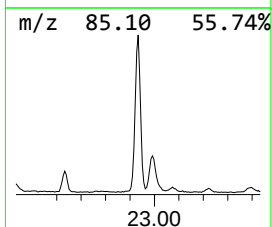
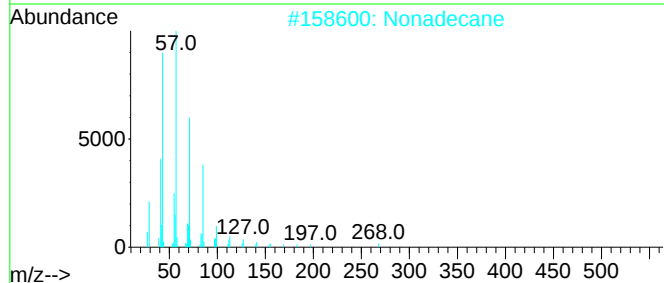
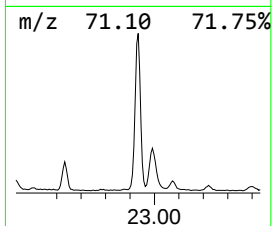
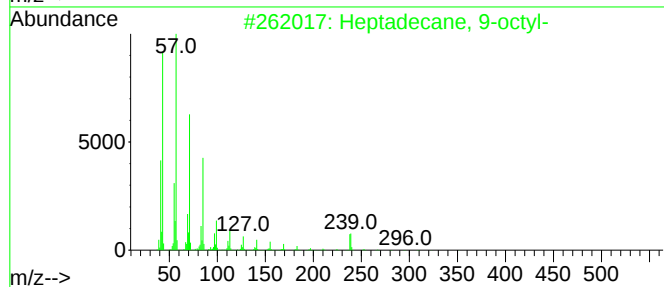
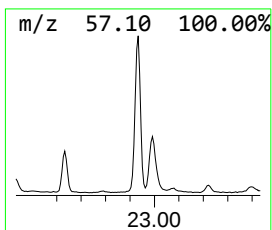
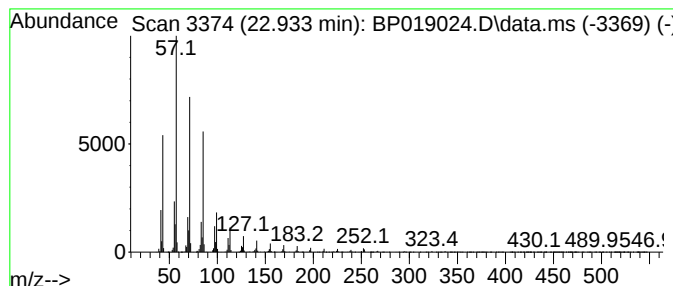
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.933	11.09 ng/ul	1523690	Perylene-d12	23.657

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
2		Nonadecane	268	C19H40	000629-92-5	94
3		Hexacosane	366	C26H54	000630-01-3	94
4		Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91
5		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019024.D
Acq On : 22 Dec 2023 05:48
Operator : MA/JU
Sample : 05808-19
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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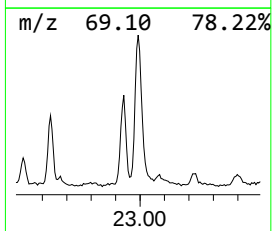
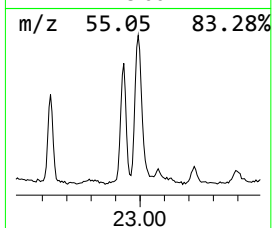
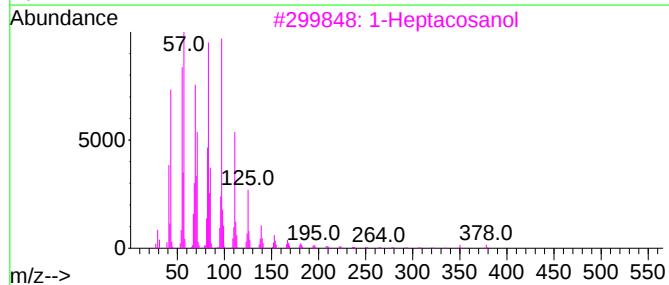
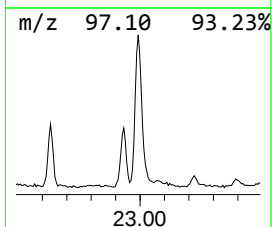
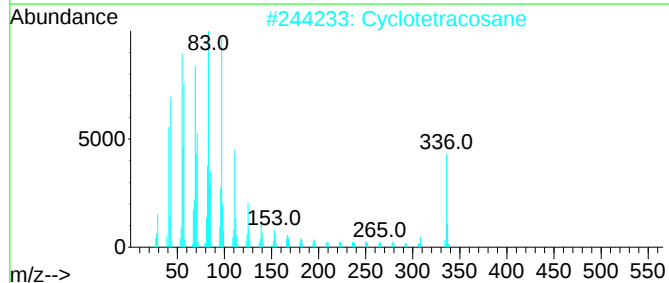
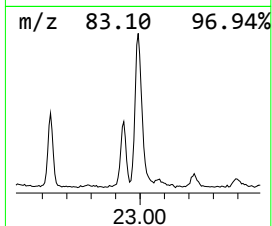
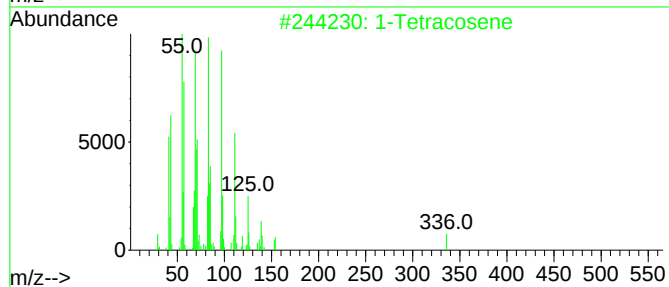
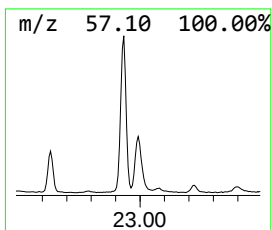
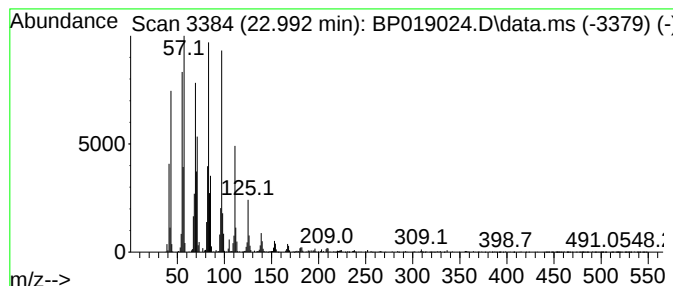
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.992	10.70 ng/ul	1470770	Perylene-d12	23.657

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Tetracosene	336	C24H48	010192-32-2	99
2		Cyclotetracosane	336	C24H48	000297-03-0	96
3		1-Heptacosanol	396	C27H56O	002004-39-9	95
4		n-Nonadecanol-1	284	C19H40O	001454-84-8	94
5		Heptacos-1-ene	378	C27H54	015306-27-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019024.D
Acq On : 22 Dec 2023 05:48
Operator : MA/JU
Sample : 05808-19
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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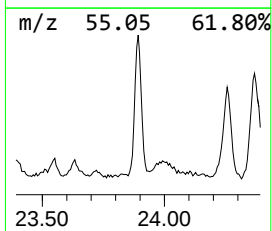
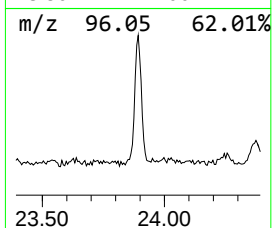
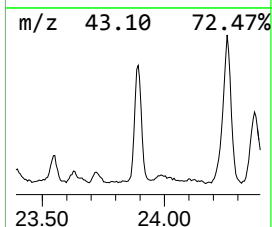
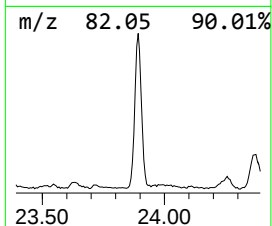
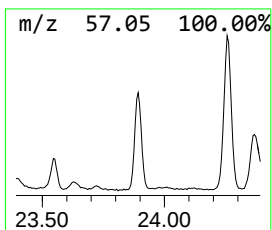
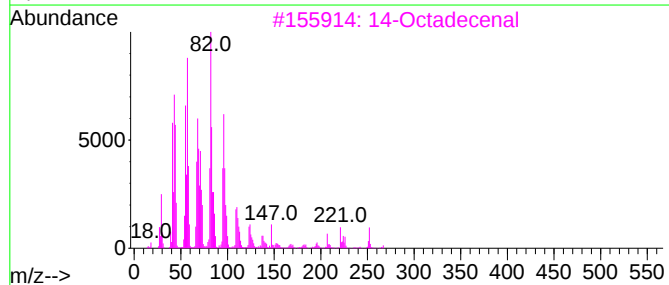
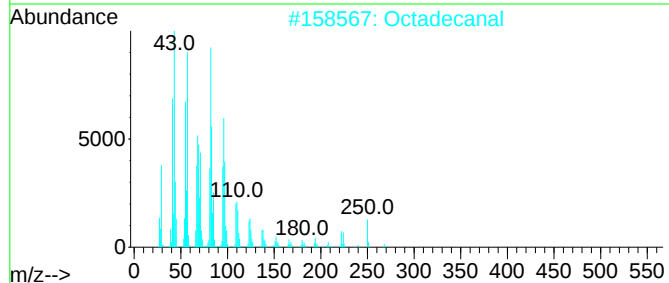
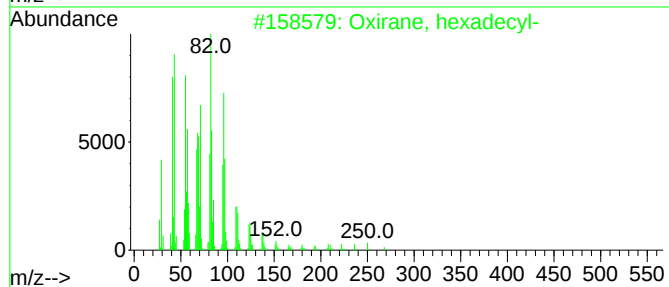
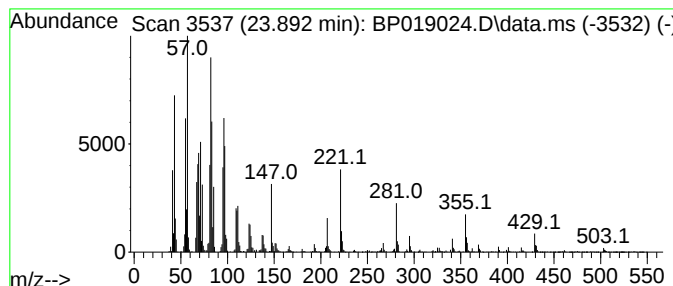
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 Oxirane, hexadecyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.892	11.90 ng/ul	1635010	Perylene-d12	23.657

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	92
2			Octadecanal	268	C18H36O	000638-66-4	91
3			14-Octadecenal	266	C18H34O	056554-89-3	86
4			1,19-Eicosadiene	278	C20H38	014811-95-1	86
5			(Z)-14-Tricosenyl formate	366	C24H46O2	077899-10-6	84



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019024.D
Acq On : 22 Dec 2023 05:48
Operator : MA/JU
Sample : 05808-19
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ClientSampleId :
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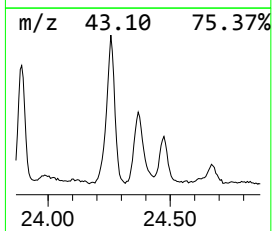
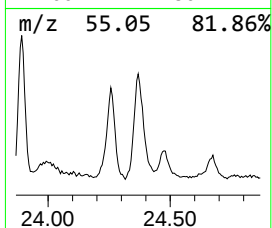
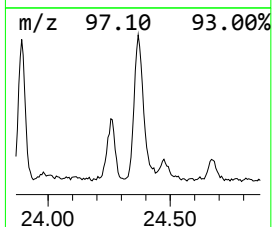
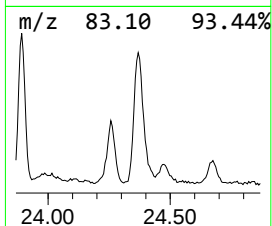
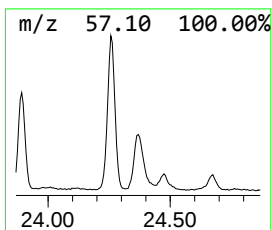
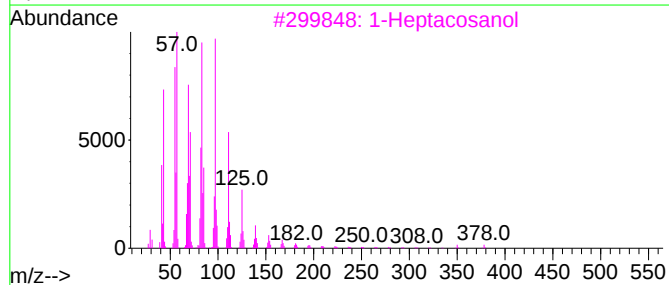
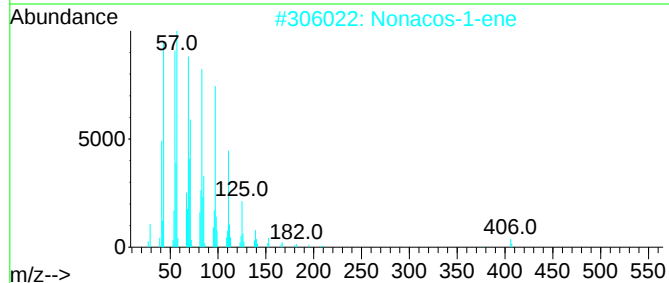
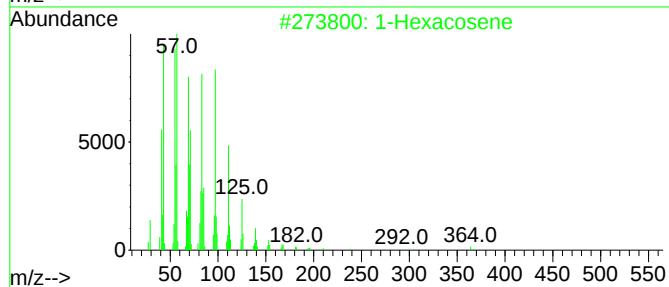
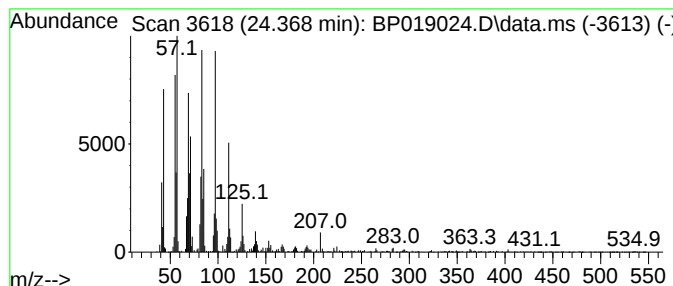
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.368	6.86 ng/ul	942666	Perylene-d12	23.657

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexacosene		364	C26H52	018835-33-1	99
2	Nonacos-1-ene		406	C29H58	018835-35-3	94
3	1-Heptacosanol		396	C27H56O	002004-39-9	94
4	9-Tricosene, (Z)-		322	C23H46	027519-02-4	93
5	Cyclotetracosane		336	C24H48	000297-03-0	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019024.D
Acq On : 22 Dec 2023 05:48
Operator : MA/JU
Sample : 05808-19
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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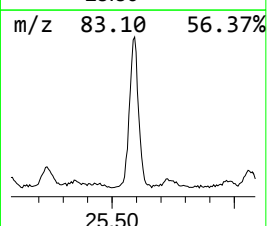
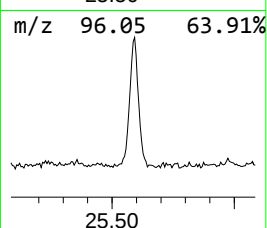
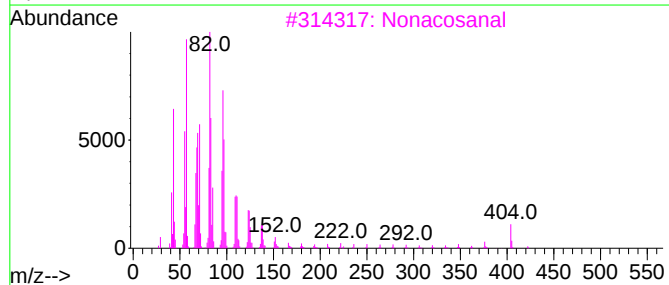
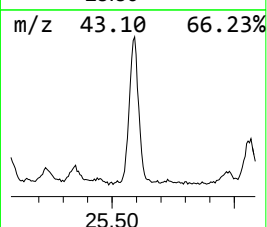
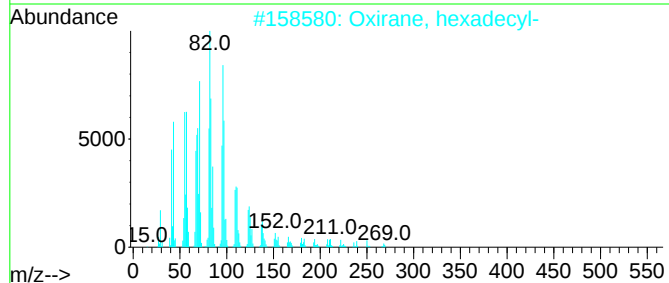
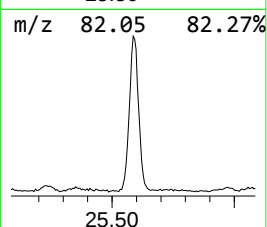
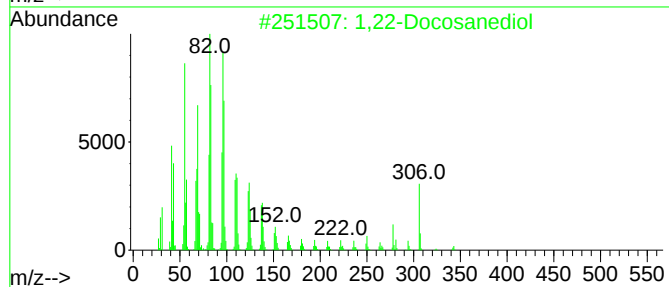
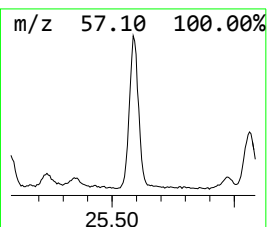
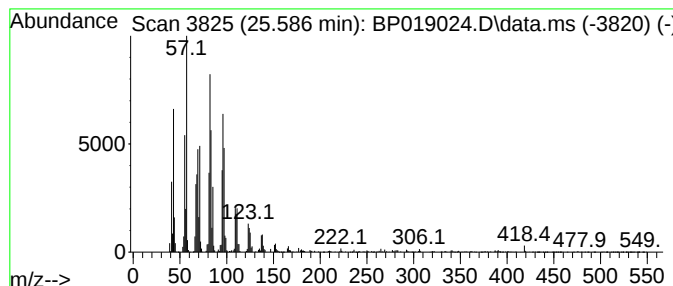
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 1,22-Docosanediol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.586	13.55 ng/ul	1862430	Perylene-d12	23.657

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,22-Docosanediol	342	C22H46O2	022513-81-1	96
2			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	94
3			Nonacosanal	422	C29H58O	072934-04-4	94
4			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
5			Triacontanal	436	C30H60O	022725-63-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019024.D
Acq On : 22 Dec 2023 05:48
Operator : MA/JU
Sample : 05808-19
Misc :
ALS Vial : 24 Sample Multiplier: 1

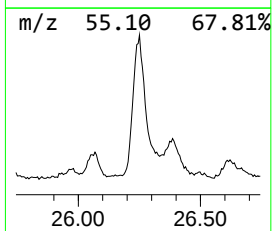
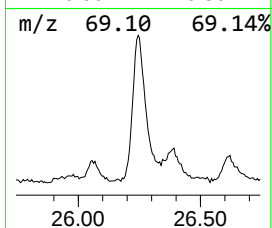
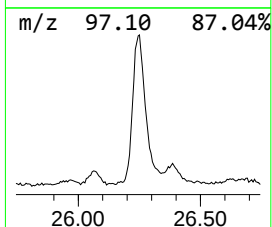
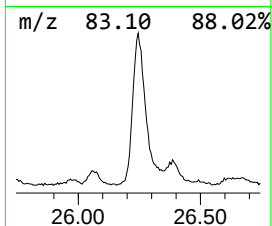
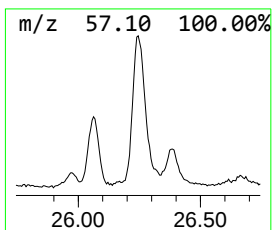
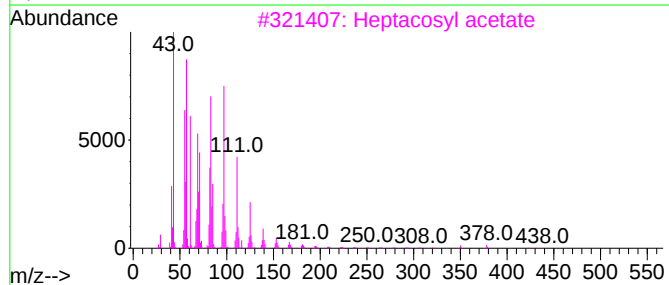
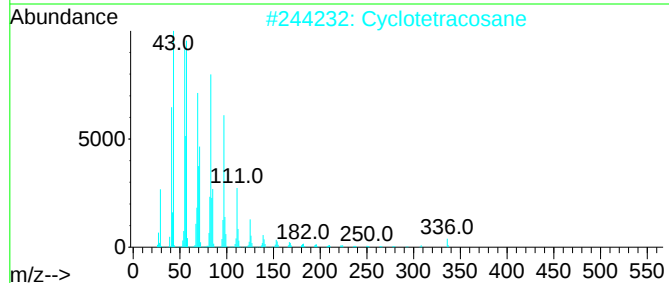
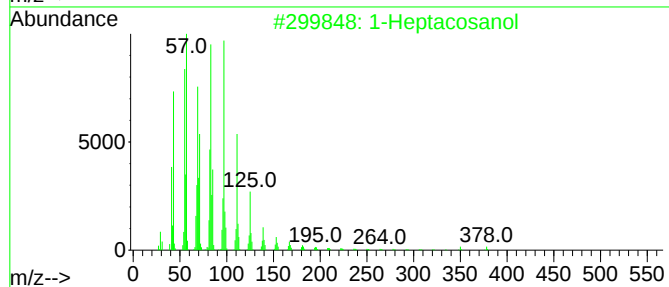
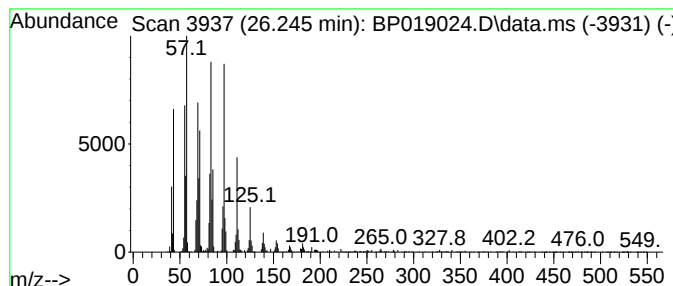
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 25 1-Heptacosanol Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
26.245	10.15 ng/ul	1394960	Perylene-d12	23.657	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heptacosanol	396	C27H56O	002004-39-9	95
2	Cyclotetracosane	336	C24H48	000297-03-0	94
3	Heptacosyl acetate	438	C29H58O2	1000351-78-2	94
4	Octacosyl acetate	452	C30H60O2	018206-97-8	93
5	1-Hexacosanol	382	C26H54O	000506-52-5	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019024.D
Acq On : 22 Dec 2023 05:48
Operator : MA/JU
Sample : 05808-19
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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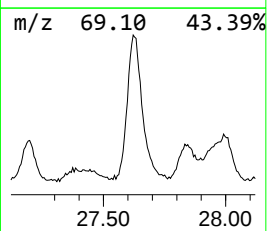
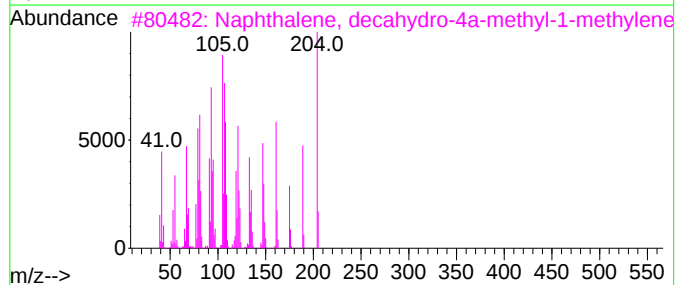
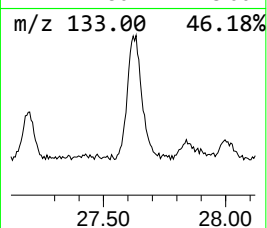
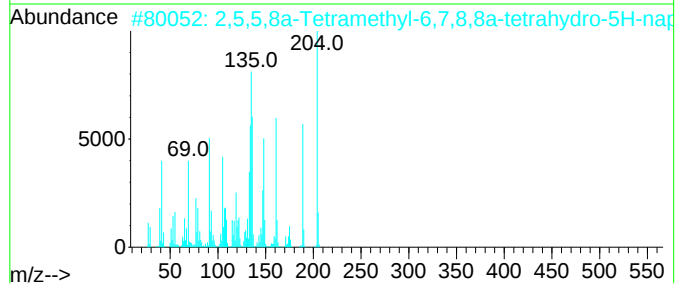
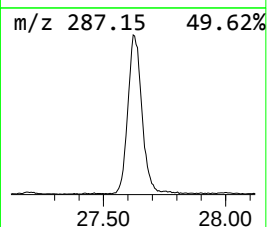
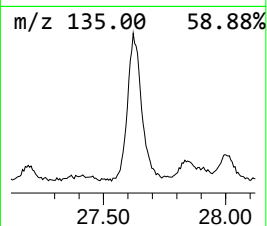
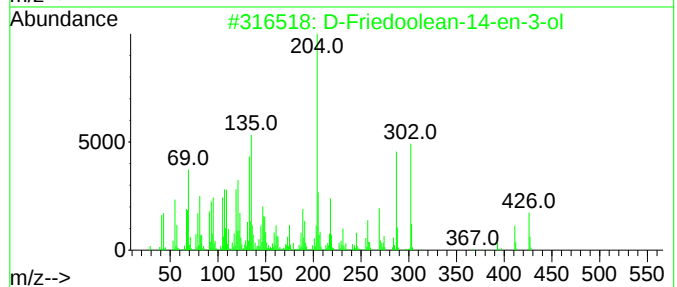
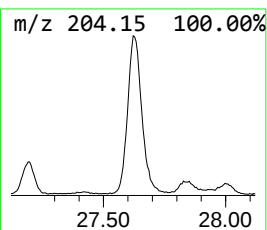
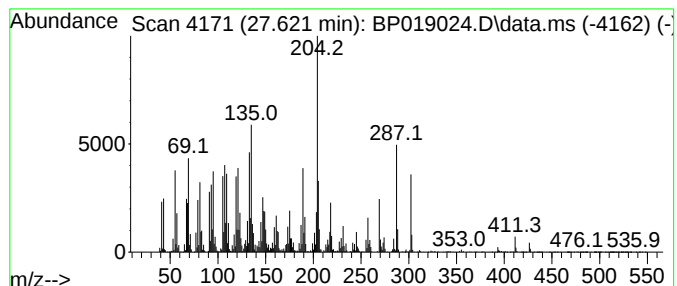
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Quant Title : SVOA CALIBRATION

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

Peak Number 26 D-Friedoolean-14-en-3-ol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.621	44.25 ng/ul	6080340	Perylene-d12	23.657

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			D-Friedoolean-14-en-3-ol	426	C30H50O	081654-73-1	74
2			2,5,5,8a-Tetramethyl-6,7,8,8a-te...	204	C14H20O	124957-09-1	49
3			Naphthalene, decahydro-4a-methyl...	204	C15H24	017066-67-0	46
4			1H-Cycloprop[e]azulene, 1a,2,3,4...	204	C15H24	000489-40-7	45
5			1,4-Dimethyl-8-isopropylidenetri...	204	C15H24	1010140-07-7	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
Data File : BP019024.D
Acq On : 22 Dec 2023 05:48
Operator : MA/JU
Sample : 05808-19
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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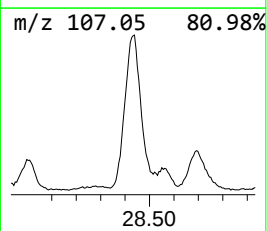
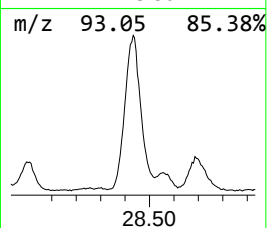
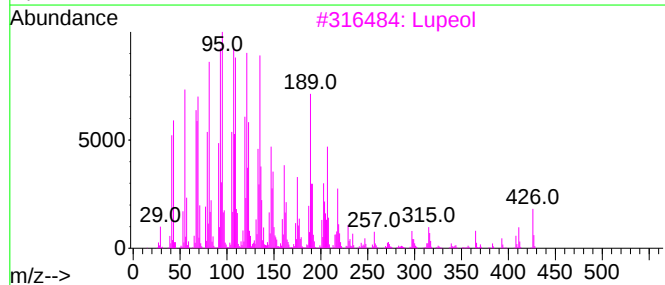
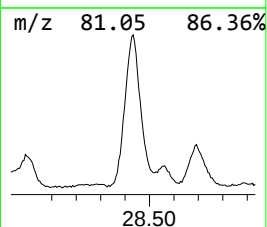
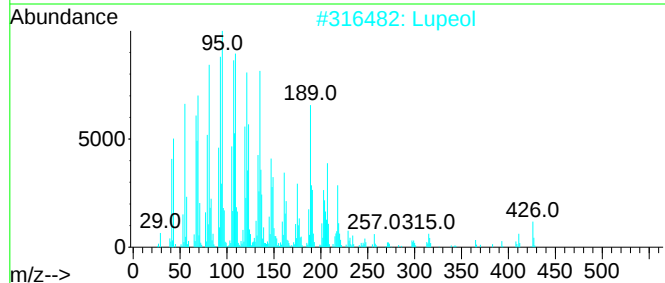
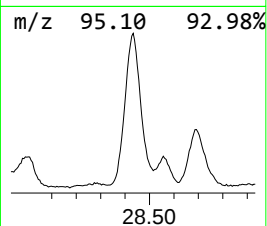
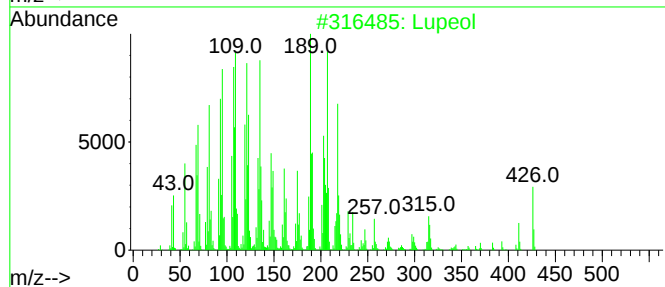
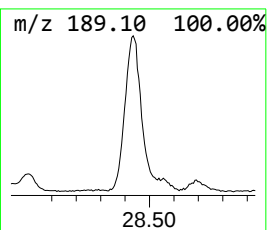
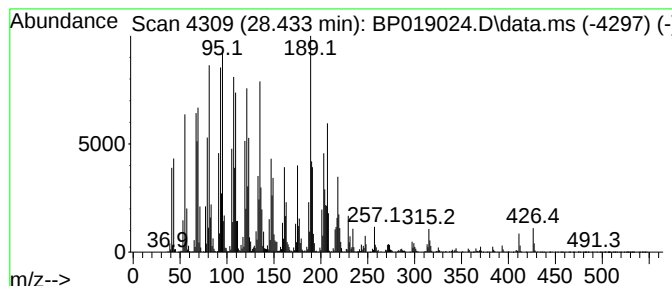
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Quant Title : SVOA CALIBRATION

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

Peak Number 27 Lupeol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.433	74.43 ng/ul	10228500	Perylene-d12	23.657

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Lupeol	426	C30H50O	000545-47-1	99
2			Lupeol	426	C30H50O	000545-47-1	95
3			Lupeol	426	C30H50O	000545-47-1	94
4			.beta.-Humulene	204	C15H24	000116-04-1	91
5			Epilupeol; 20(29)-Lupen-3alpha-o...	426	C30H50O	1000513-01-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
 Data File : BP019024.D
 Acq On : 22 Dec 2023 05:48
 Operator : MA/JU
 Sample : 05808-19
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.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
Benzeneacetic a...	8.763	3.7	ng/ul	161325	1	7.816	868160	20.0
Benzoic acid	9.916	12.0	ng/ul	712552	2	10.610	1190740	20.0
2-Phenylpropenal	10.063	3.7	ng/ul	220493	2	10.610	1190740	20.0
unknown-01	11.169	7.1	ng/ul	423241	2	10.610	1190740	20.0
N-Methoxy-N-met...	11.404	2.8	ng/ul	164649	2	10.610	1190740	20.0
n-Hexadecanoic ...	18.098	6.3	ng/ul	679229	4	17.204	2139730	20.0
N-(2-Cyanopheny...	18.651	4.4	ng/ul	473494	4	17.204	2139730	20.0
.beta.-Phenylpr...	18.869	4.0	ng/ul	431238	4	17.204	2139730	20.0
Anthracene, 9-p...	18.916	2.1	ng/ul	219938	4	17.204	2139730	20.0
Dibenzoylmethane	19.122	16.7	ng/ul	1790210	4	17.204	2139730	20.0
Octadecanoic acid	19.333	2.0	ng/ul	324457	5	21.298	3167880	20.0
E-14-Hexadecenal	20.051	2.8	ng/ul	440838	5	21.298	3167880	20.0
5-Bromo-4-oxo-4...	20.833	7.3	ng/ul	1153050	5	21.298	3167880	20.0
(DEL) Alkane: S...	21.016	16.2	ng/ul	2566950	5	21.298	3167880	20.0
(DEL) Alkane: S...	21.922	41.2	ng/ul	6522120	5	21.298	3167880	20.0
Tetracosanoic acid	22.251	5.0	ng/ul	787338	5	21.298	3167880	20.0
unknown-02	22.433	3.4	ng/ul	536274	5	21.298	3167880	20.0
Hexacosanal	22.633	6.1	ng/ul	842991	6	23.657	2748370	20.0
1,3,5-Triphenyl...	22.822	3.0	ng/ul	410571	6	23.657	2748370	20.0
(DEL) Alkane: S...	22.933	11.1	ng/ul	1523690	6	23.657	2748370	20.0
(DEL) Alkane: S...	22.992	10.7	ng/ul	1470770	6	23.657	2748370	20.0
Oxirane, hexade...	23.892	11.9	ng/ul	1635010	6	23.657	2748370	20.0
(DEL) Alkane: S...	24.368	6.9	ng/ul	942666	6	23.657	2748370	20.0
1,22-Docosanediol	25.586	13.6	ng/ul	1862430	6	23.657	2748370	20.0
1-Heptacosanol	26.245	10.2	ng/ul	1394960	6	23.657	2748370	20.0
D-Friedoolean-1...	27.621	44.3	ng/ul	6080340	6	23.657	2748370	20.0
Lupeol	28.433	74.4	ng/ul	10228500	6	23.657	2748370	20.0