

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
 Data File : BP018124.D
 Acq On : 13 Nov 2023 02:47
 Operator : MA/JU
 Sample : 05158-02
 Misc :
 ALS Vial : 105 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH363

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-BP111023.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP018124.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.228	20	24	34	rVB3	16960	25869	1.63%	0.123%
2	3.669	96	99	114	rBV	73961	137851	8.67%	0.654%
3	3.963	146	149	152	rVB3	7262	8334	0.52%	0.040%
4	4.481	233	237	240	rBV6	4815	8174	0.51%	0.039%
5	4.511	240	242	246	rVB4	2890	3699	0.23%	0.018%
6	4.752	280	283	285	rBV4	2735	2748	0.17%	0.013%
7	5.387	387	391	397	rVB3	13611	19617	1.23%	0.093%
8	6.593	591	596	597	rBV4	3312	5669	0.36%	0.027%
9	7.022	663	669	687	rBV2	42151	116670	7.33%	0.554%
10	7.193	692	698	710	rBV	217726	357639	22.48%	1.697%
11	7.363	719	727	744	rBV	193821	362422	22.78%	1.720%
12	7.687	777	782	784	rBV5	1893	2821	0.18%	0.013%
13	7.840	801	808	822	rBV	419746	693084	43.57%	3.289%
14	8.187	863	867	869	rBV5	2173	2455	0.15%	0.012%
15	8.504	918	921	924	rBV5	1854	2403	0.15%	0.011%
16	8.575	927	933	948	rBV	127586	291788	18.34%	1.385%
17	8.716	955	957	959	rVB3	2317	2197	0.14%	0.010%
18	8.816	971	974	977	rBV4	1422	2566	0.16%	0.012%
19	9.051	1007	1014	1028	rBV	259149	492200	30.94%	2.336%
20	9.216	1037	1042	1050	rVV	55935	85101	5.35%	0.404%
21	9.269	1050	1051	1055	rVB4	2300	2445	0.15%	0.012%
22	9.393	1068	1072	1077	rBV8	3110	5990	0.38%	0.028%
23	9.598	1104	1107	1113	rBV8	1791	2712	0.17%	0.013%
24	9.775	1132	1137	1149	rBV2	36458	79732	5.01%	0.378%
25	9.875	1153	1154	1158	rVV4	2141	2370	0.15%	0.011%
26	9.934	1158	1164	1169	rVB8	4843	9019	0.57%	0.043%
27	10.298	1220	1226	1242	rBV	129728	311615	19.59%	1.479%
28	10.657	1279	1287	1302	rBV	569163	982137	61.74%	4.661%
29	10.845	1312	1319	1340	rBV	247085	567843	35.70%	2.695%
30	11.140	1364	1369	1372	rBV6	1967	3875	0.24%	0.018%
31	11.492	1427	1429	1430	rBV2	2658	2174	0.14%	0.010%
32	11.534	1433	1436	1439	rVV5	1825	2238	0.14%	0.011%
33	11.704	1460	1465	1473	rVV8	8393	17494	1.10%	0.083%
34	11.957	1504	1508	1509	rBV3	1953	2371	0.15%	0.011%
35	11.975	1509	1511	1514	rVV3	2060	2595	0.16%	0.012%
36	12.269	1556	1561	1567	rBV7	4061	8150	0.51%	0.039%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
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37	12.334	1567	1572	1582	rBV4	17130	51430	3.23%	0.244%
38	12.704	1632	1635	1639	rVB6	1625	2511	0.16%	0.012%
39	12.751	1639	1643	1646	rBV6	4185	5379	0.34%	0.026%
40	13.257	1726	1729	1731	rVV4	2600	2772	0.17%	0.013%

41	13.357	1740	1746	1754	rBV3	18933	34435	2.16%	0.163%
42	13.545	1774	1778	1782	rVB7	2200	3379	0.21%	0.016%
43	13.622	1787	1791	1794	rVB5	2174	3116	0.20%	0.015%
44	13.722	1801	1808	1813	rBV5	3852	6325	0.40%	0.030%
45	13.816	1820	1824	1826	rBV5	1733	2312	0.15%	0.011%

46	13.851	1826	1830	1831	rBV4	2754	3307	0.21%	0.016%
47	13.939	1839	1845	1852	rBV	673672	961848	60.46%	4.565%
48	14.134	1873	1878	1882	rBV8	2828	5205	0.33%	0.025%
49	14.216	1885	1892	1906	rBV	724767	1052347	66.15%	4.994%
50	14.434	1927	1929	1934	rVB5	2291	3048	0.19%	0.014%

51	14.516	1934	1943	1955	rBV	884130	1259313	79.16%	5.976%
52	15.028	2027	2030	2035	rBV7	3352	4945	0.31%	0.023%
53	15.133	2044	2048	2051	rBV6	2698	3597	0.23%	0.017%
54	15.263	2066	2070	2072	rBV5	2042	2806	0.18%	0.013%
55	15.298	2072	2076	2081	rVB8	4359	6263	0.39%	0.030%

56	15.516	2105	2113	2126	rBV	864812	1313701	82.58%	6.234%
57	15.757	2153	2154	2162	rVB8	4721	7103	0.45%	0.034%
58	15.816	2162	2164	2168	rVB5	2653	4045	0.25%	0.019%
59	15.886	2173	2176	2181	rVB7	2974	4090	0.26%	0.019%
60	16.169	2220	2224	2227	rBV5	4311	4994	0.31%	0.024%

61	16.257	2235	2239	2242	rBV6	3735	6951	0.44%	0.033%
62	16.322	2248	2250	2254	rVB5	1781	2149	0.14%	0.010%
63	16.663	2304	2308	2310	rBV4	4286	5183	0.33%	0.025%
64	16.851	2336	2340	2347	rBV9	3996	9383	0.59%	0.045%
65	16.975	2358	2361	2365	rVB6	4479	5562	0.35%	0.026%

66	17.086	2377	2380	2384	rBV6	5111	7428	0.47%	0.035%
67	17.163	2389	2393	2395	rBV5	2691	4266	0.27%	0.020%
68	17.257	2405	2409	2411	rBV4	22604	28944	1.82%	0.137%
69	17.298	2411	2416	2425	rVV	869310	1296558	81.50%	6.153%
70	17.398	2427	2433	2448	rVB	901242	1354835	85.17%	6.430%

71	17.580	2462	2464	2470	rVB7	7927	10996	0.69%	0.052%
72	17.710	2483	2486	2490	rBV6	4357	7332	0.46%	0.035%
73	17.916	2519	2521	2528	rVB8	6414	9886	0.62%	0.047%
74	18.145	2554	2560	2565	rBV8	19104	33680	2.12%	0.160%
75	18.239	2572	2576	2580	rVB	18770	22780	1.43%	0.108%

76	18.304	2585	2587	2588	rBV2	2455	2374	0.15%	0.011%
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77	18.351	2593	2595	2597	rBV3	2825	2435	0.15%	0.012%
78	18.386	2599	2601	2602	rBV2	2453	2256	0.14%	0.011%
79	18.527	2622	2625	2627	rBV4	4004	4211	0.26%	0.020%
80	18.563	2627	2631	2637	rVB8	18748	27134	1.71%	0.129%
81	18.633	2638	2643	2648	rBV	154046	188055	11.82%	0.892%
82	18.769	2664	2666	2670	rVB5	5024	5673	0.36%	0.027%
83	18.810	2670	2673	2676	rBV4	6330	7040	0.44%	0.033%
84	18.904	2687	2689	2691	rBV3	3433	3215	0.20%	0.015%
85	19.127	2723	2727	2732	rVB6	7679	11946	0.75%	0.057%
86	19.227	2739	2744	2748	rBV6	15409	24428	1.54%	0.116%
87	19.298	2753	2756	2760	rVB2	8453	13031	0.82%	0.062%
88	19.521	2791	2794	2797	rBV5	8220	10124	0.64%	0.048%
89	19.651	2811	2816	2836	rBV	1157458	1514055	95.18%	7.185%
90	19.945	2864	2866	2872	rVB6	11172	9159	0.58%	0.043%
91	21.086	3055	3060	3068	rBV	153095	260158	16.35%	1.235%
92	21.427	3113	3118	3127	rBV	877627	1205229	75.76%	5.720%
93	22.163	3240	3243	3249	rVB	104136	138470	8.70%	0.657%
94	22.292	3260	3265	3271	rBV	683776	886250	55.71%	4.206%
95	22.568	3308	3312	3320	rVB2	240834	356825	22.43%	1.693%
96	22.815	3351	3354	3361	rVB4	89793	140760	8.85%	0.668%
97	23.098	3397	3402	3408	rVB	506402	770152	48.41%	3.655%
98	23.292	3430	3435	3442	rVB4	207768	368329	23.15%	1.748%
99	23.757	3508	3514	3523	rVB2	756810	1590799	100.00%	7.550%
100	23.927	3536	3543	3554	rVB	604790	1349108	84.81%	6.403%

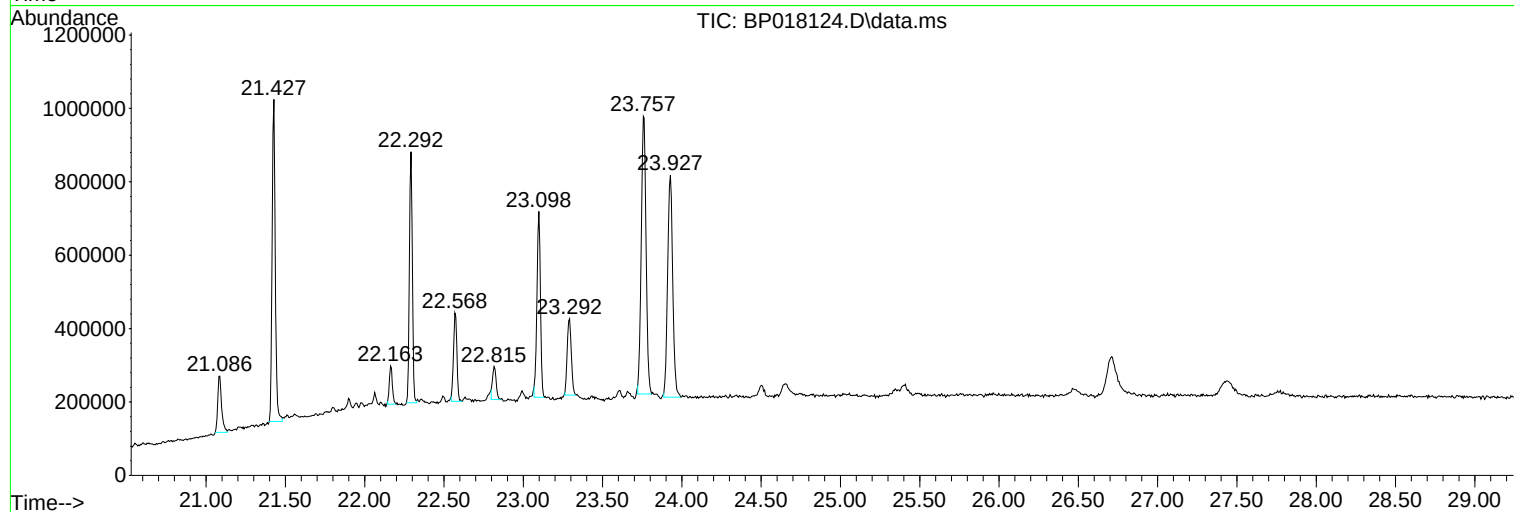
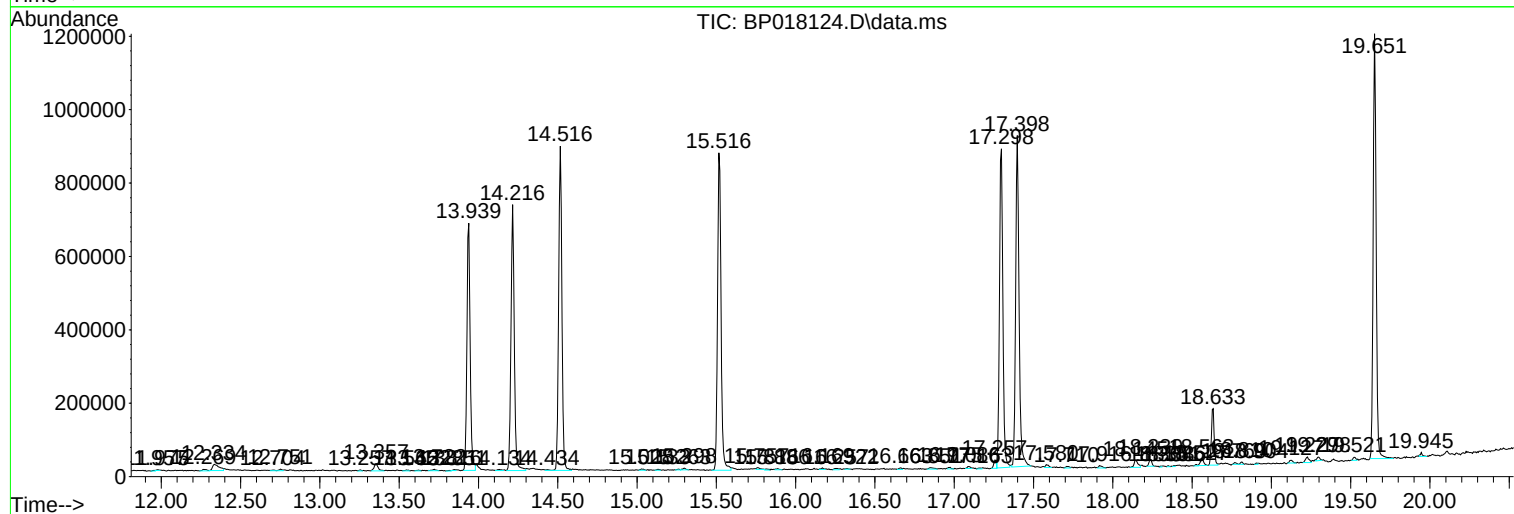
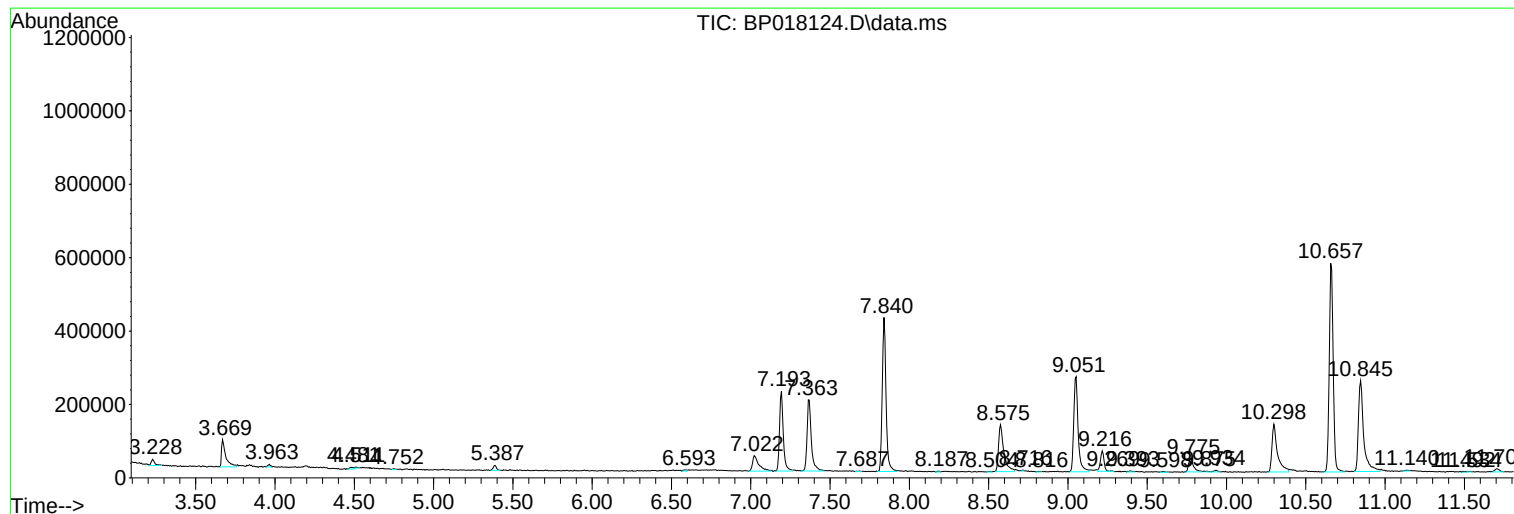
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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



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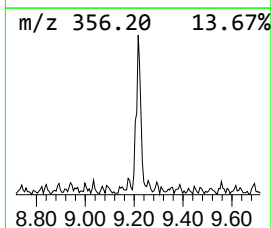
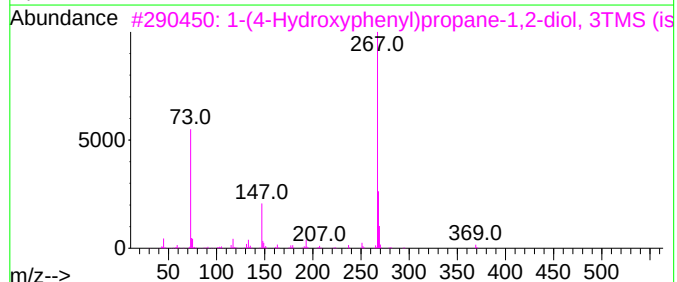
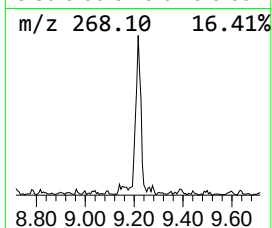
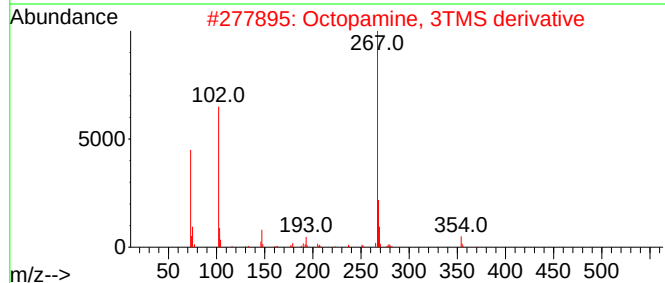
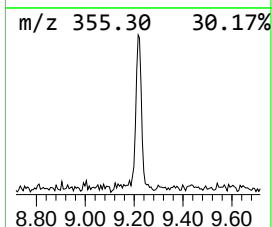
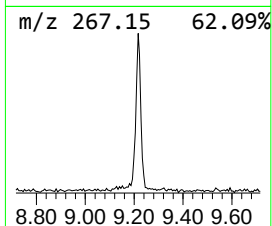
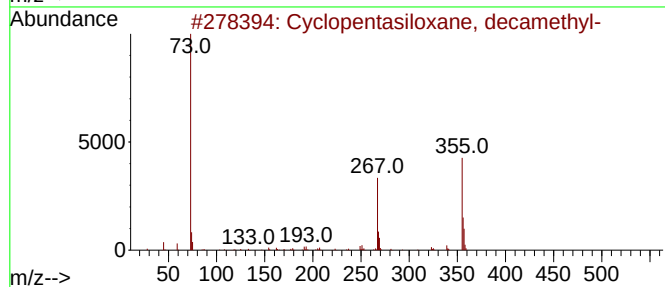
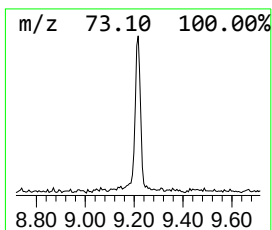
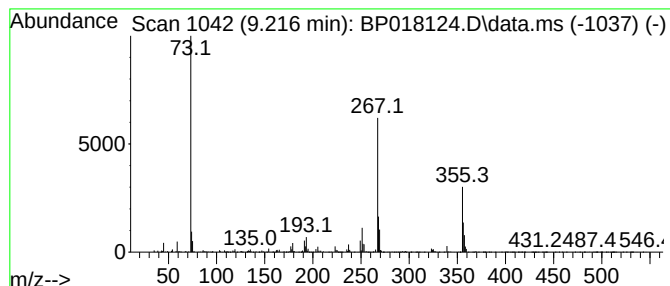
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Cyclopentasiloxane, decamet... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.216	2.46 ng/ul	85101	1,4-Dichlorobenzene-d4	7.840

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	86
2			Octopamine, 3TMS derivative	369	C17H35NO2Si3	068595-60-8	50
3			1-(4-Hydroxyphenyl)propane-1,2-d...	384	C18H36O3Si3	1000495-85-7	47
4			[(4-Hexylbenzene-1,3-diyl)bis(ox...	338	C18H34O2Si2	134273-01-1	47
5			2-Amino-4-methylphenol, 2TMS der...	267	C13H25NOSi2	1000373-28-1	47



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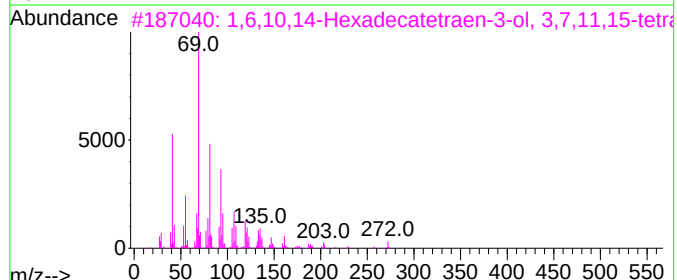
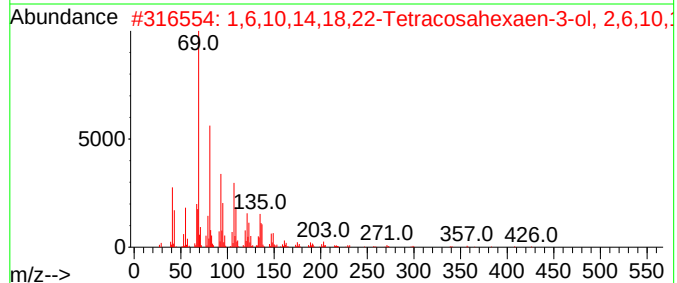
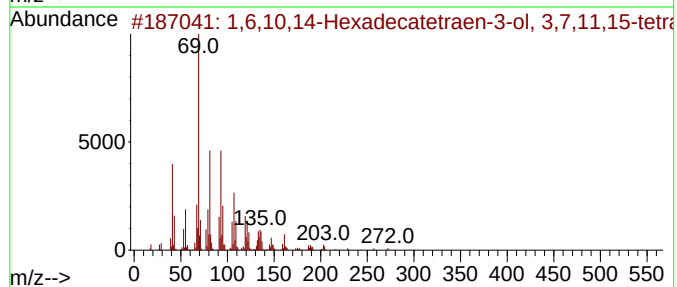
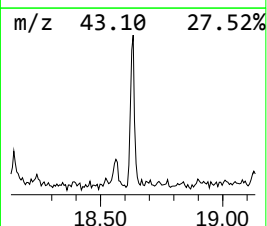
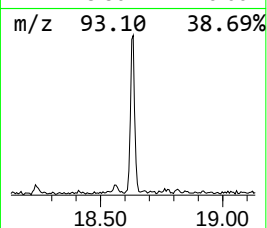
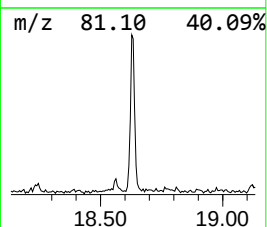
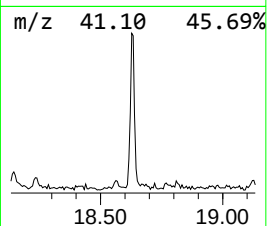
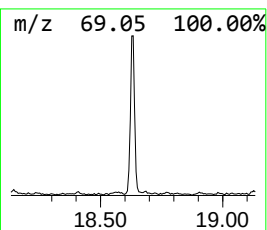
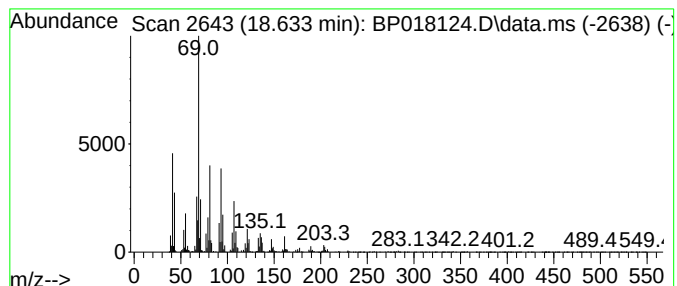
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1,6,10,14-Hexadecatetraen-3... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.633	2.90 ng/ul	188055	Phenanthrene-d10	17.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,6,10,14-Hexadecatetraen-3-ol, ...	290	C20H34O	001113-21-9	83		
2	1,6,10,14,18,22-Tetracosahexaen-...	426	C30H50O	097232-74-1	74		
3	1,6,10,14-Hexadecatetraen-3-ol, ...	290	C20H34O	001113-21-9	64		
4	1,6,10-Dodecatrien-3-ol, 3,7,11-...	222	C15H26O	007212-44-4	64		
5	(2E,6E,10E)-3,7,11,15-Tetramethy...	318	C21H34O2	125456-63-5	64		



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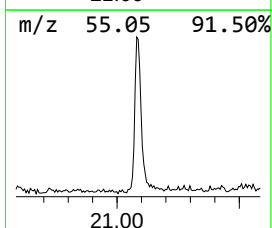
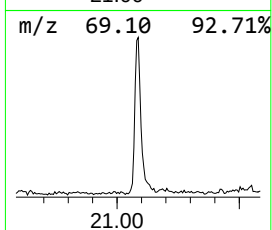
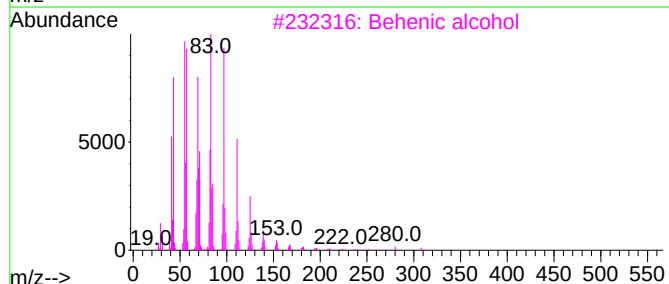
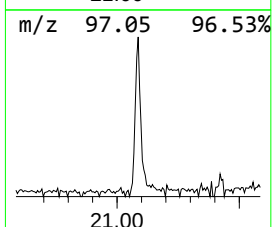
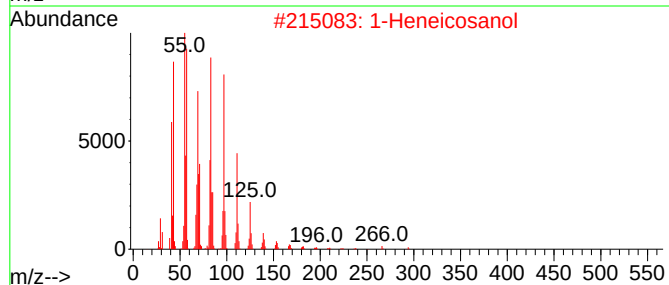
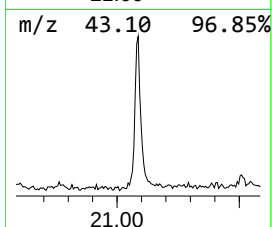
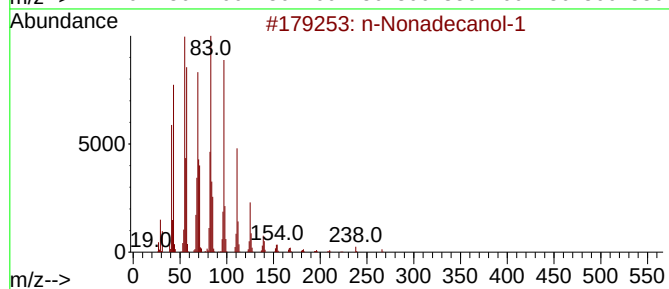
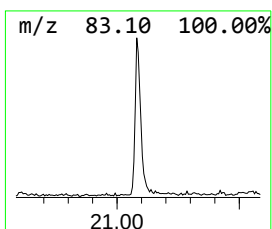
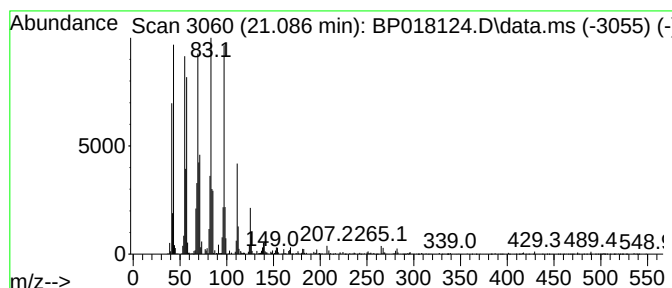
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Nonadecanol-1 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.086	4.32 ng/ul	260158	Chrysene-d12	21.427	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Nonadecanol-1	284	C19H40O	001454-84-8	94
2	1-Heneicosanol	312	C21H44O	015594-90-8	94
3	Behenic alcohol	326	C22H46O	000661-19-8	94
4	1-Tricosene	322	C23H46	018835-32-0	94
5	n-Nonadecanol-1	284	C19H40O	001454-84-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018124.D
Acq On : 13 Nov 2023 02:47
Operator : MA/JU
Sample : 05158-02
Misc :
ALS Vial : 105 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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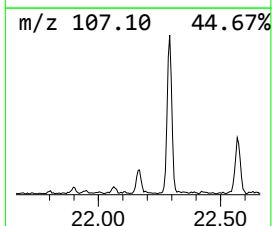
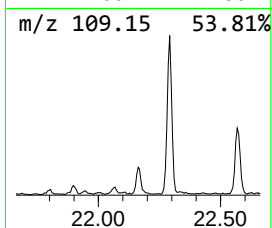
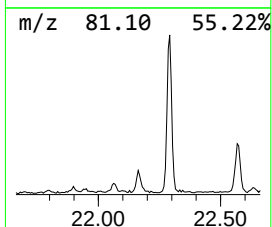
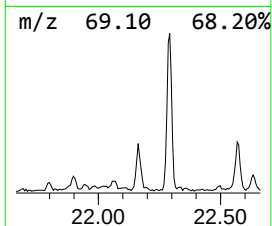
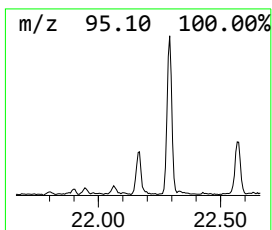
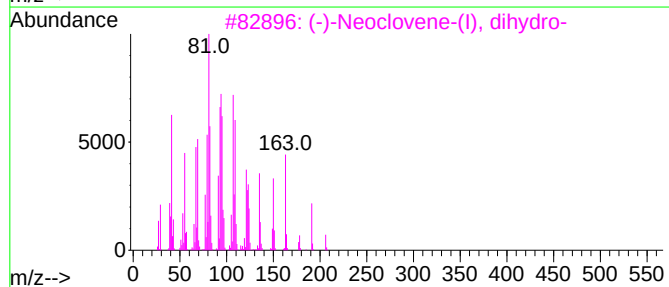
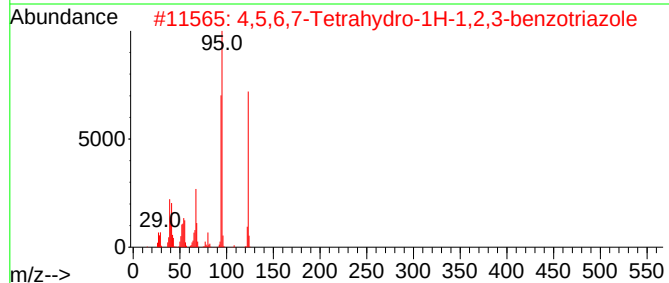
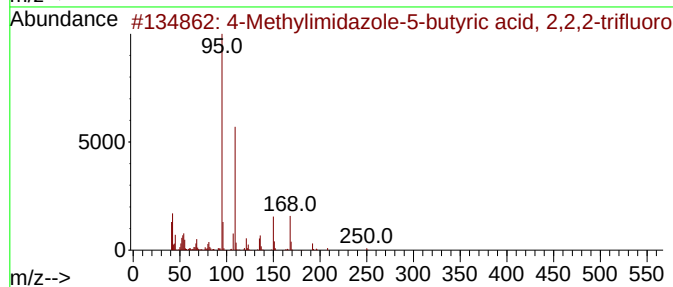
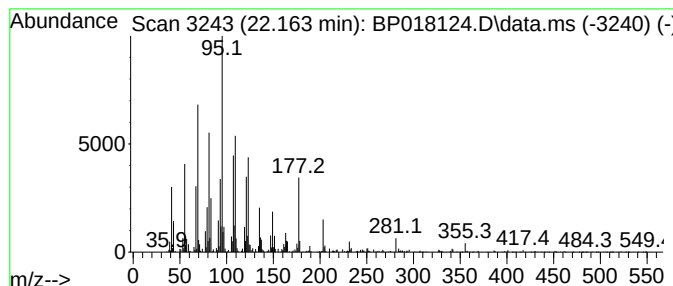
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.163	2.30 ng/ul	138470	Chrysene-d12	21.427

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Methylimidazole-5-butyric acid...	250	C10H13F3N2O2	1000129-15-9	38
2			4,5,6,7-Tetrahydro-1H-1,2,3-benz...	123	C6H9N3	006789-99-7	35
3			(-)-Neoclovene-(I), dihydro-	206	C15H26	1000152-82-1	27
4			Marrubiin	332	C20H28O4	000465-92-9	27
5			4,7-Methanobenzofuran, 2,2'-oxyb...	374	C24H38O3	081955-10-4	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018124.D
Acq On : 13 Nov 2023 02:47
Operator : MA/JU
Sample : 05158-02
Misc :
ALS Vial : 105 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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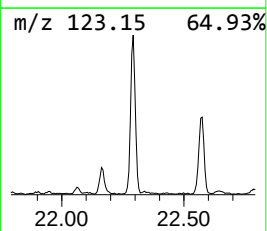
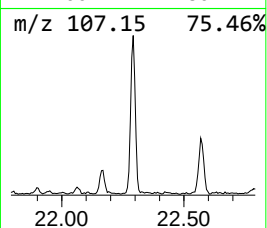
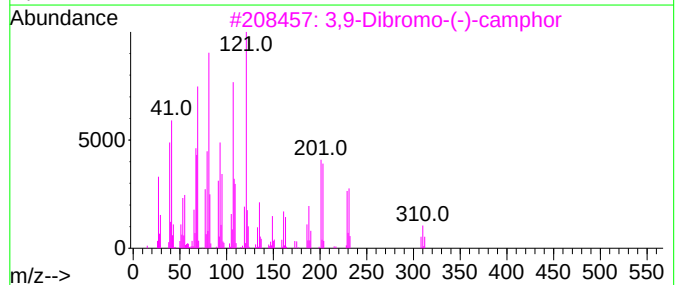
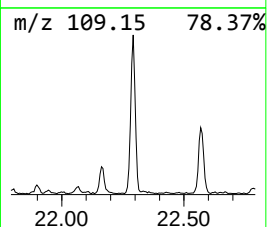
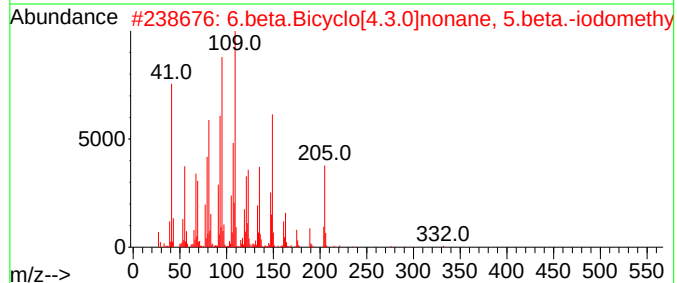
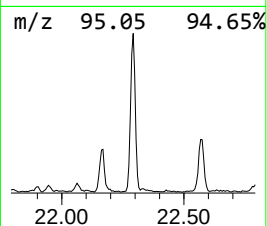
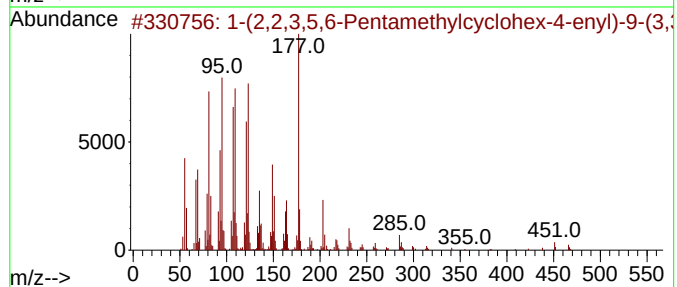
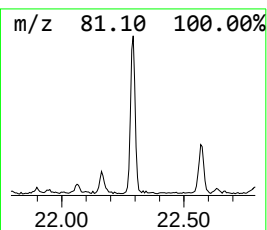
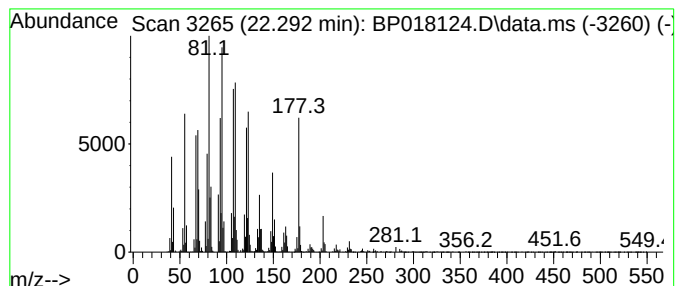
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1-(2,2,3,5,6-Pentamethylcyc... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.292	14.71 ng/ul	886250	Chrysene-d12	21.427

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-(2,2,3,5,6-Pentamethylcyclohex...	466	C34H58	1000373-94-7	90		
2	6.beta.Bicyclo[4.3.0]nonane, 5.b...	332	C15H25I	1000195-85-9	49		
3	3,9-Dibromo-(-)-camphor	308	C10H14Br2O	064474-55-1	38		
4	Cyclodecene, 3-bromo-	216	C10H17Br	056325-56-5	25		
5	1-Isopropenyl-3-propenylcyclopent...	150	C11H18	1000192-81-2	22		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018124.D
Acq On : 13 Nov 2023 02:47
Operator : MA/JU
Sample : 05158-02
Misc :
ALS Vial : 105 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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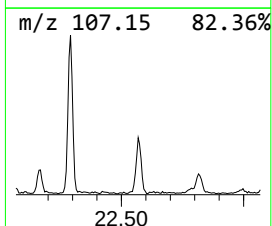
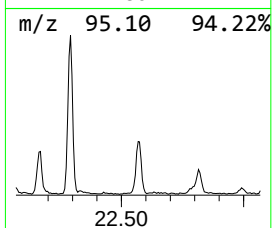
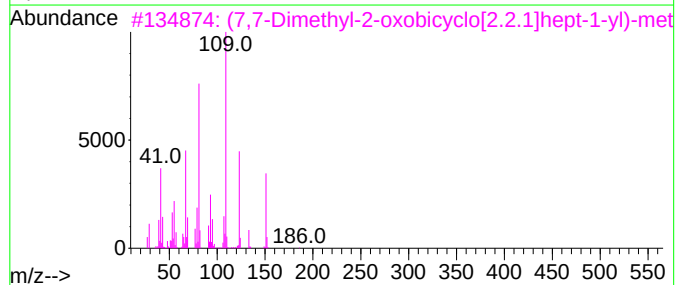
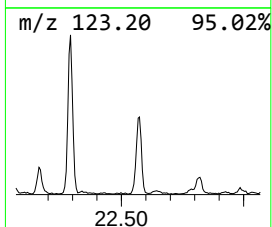
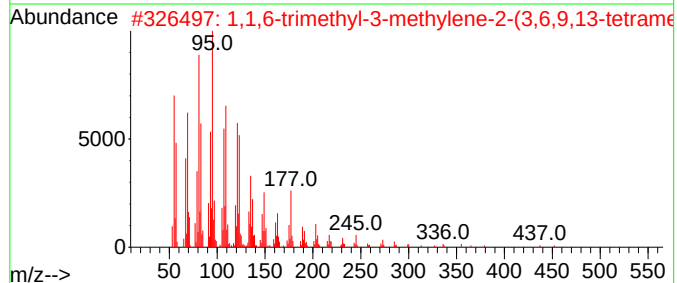
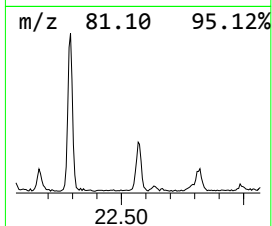
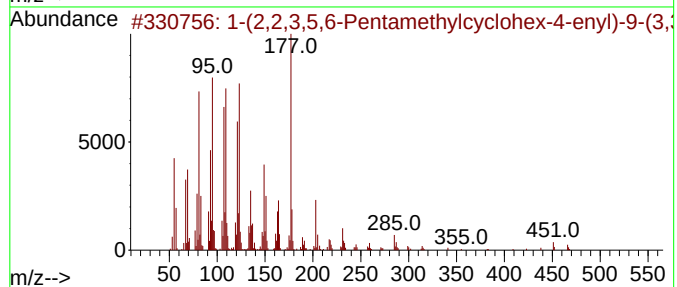
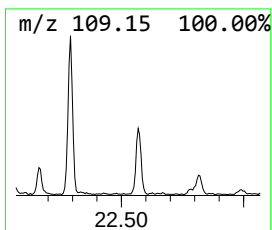
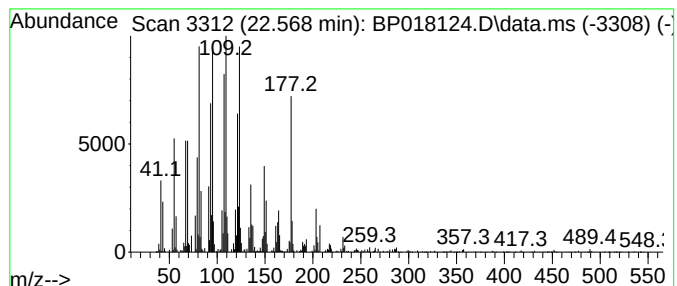
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1,1,6-trimethyl-3-methylene... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.568	5.92 ng/ul	356825	Chrysene-d12	21.427

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-(2,2,3,5,6-Pentamethylcyclohex...	466	C34H58	1000373-94-7	91
2			1,1,6-trimethyl-3-methylene-2-(3...	452	C33H56	1000373-94-5	53
3			(7,7-Dimethyl-2-oxobicyclo[2.2.1...	250	C10H15ClO3S	1000194-76-1	27
4			Diethyl allylphosphonate	178	C7H15O3P	001067-87-4	25
5			Cyclohexane, 1,2-dimethyl-3,5-bi...	192	C14H24	074806-55-6	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018124.D
Acq On : 13 Nov 2023 02:47
Operator : MA/JU
Sample : 05158-02
Misc :
ALS Vial : 105 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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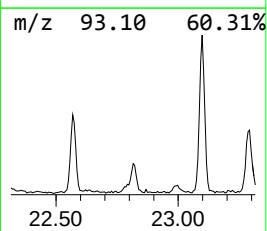
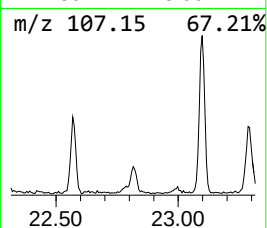
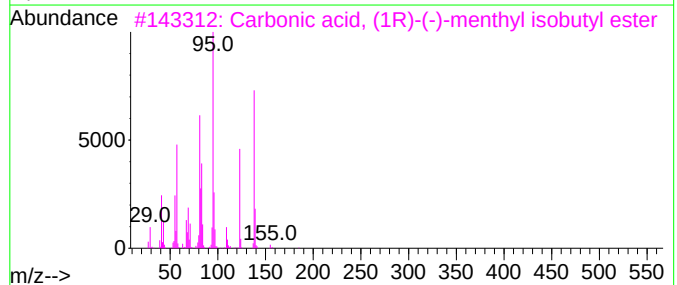
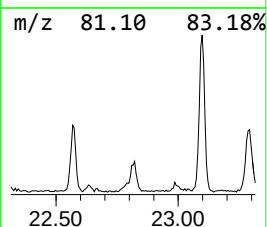
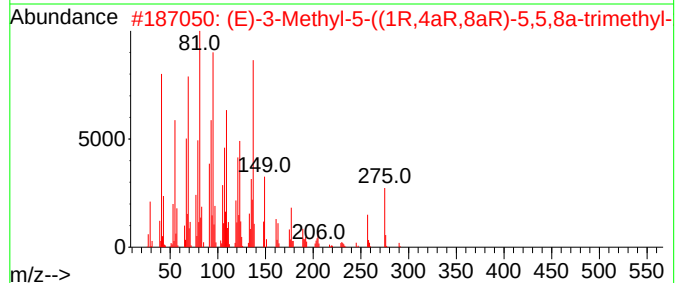
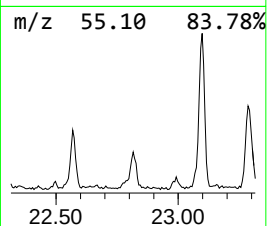
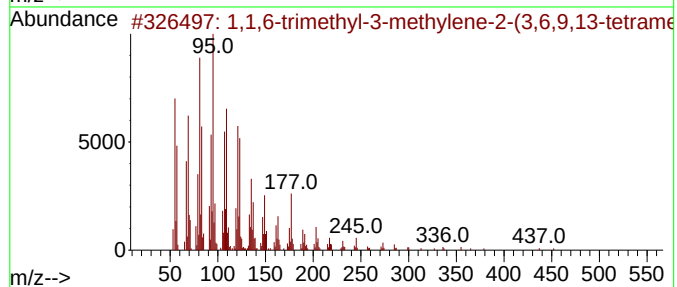
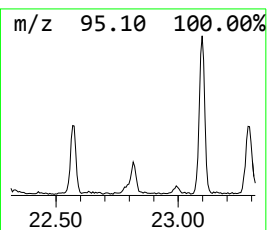
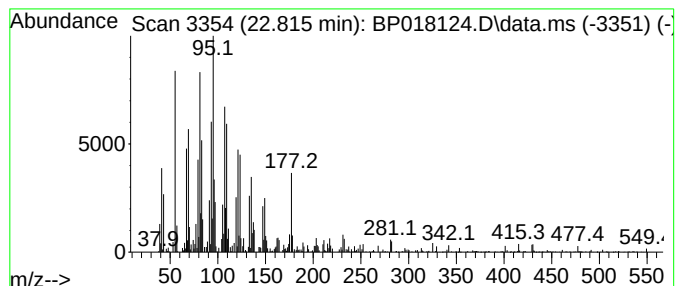
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Carbonic acid, (1R)-(-)-men... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.815	2.09 ng/ul	140760	Perylene-d12	23.927

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1,6-trimethyl-3-methylene-2-(3...	452	C33H56	1000373-94-5	78		
2	(E)-3-Methyl-5-((1R,4aR,8aR)-5,5...	290	C20H34O	021738-29-4	53		
3	Carbonic acid, (1R)-(-)-menthyl ...	256	C15H28O3	1010392-44-8	53		
4	Ergost-25-ene-3,5,6,12-tetrol, (...)	448	C28H48O4	056052-97-2	49		
5	Divinylbis(cyclopropyl)silane	164	C10H16Si	1000109-95-2	38		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018124.D
Acq On : 13 Nov 2023 02:47
Operator : MA/JU
Sample : 05158-02
Misc :
ALS Vial : 105 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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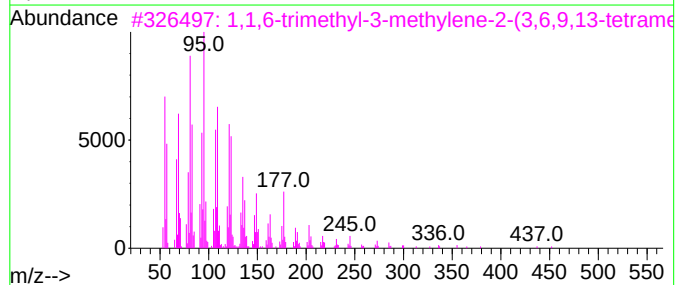
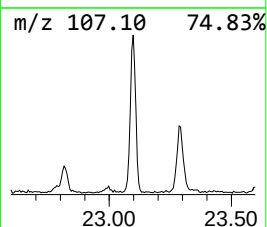
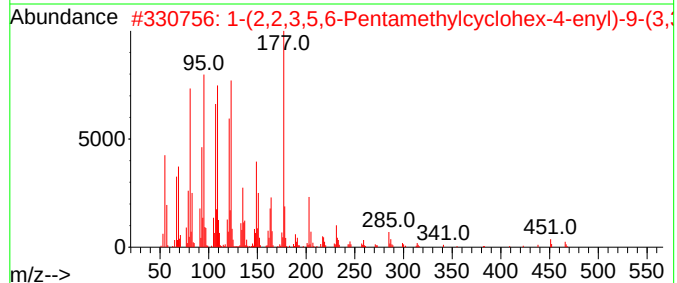
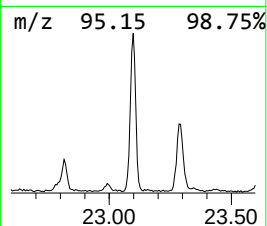
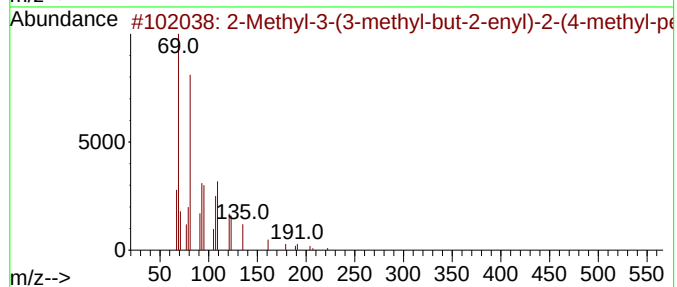
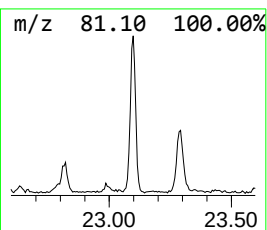
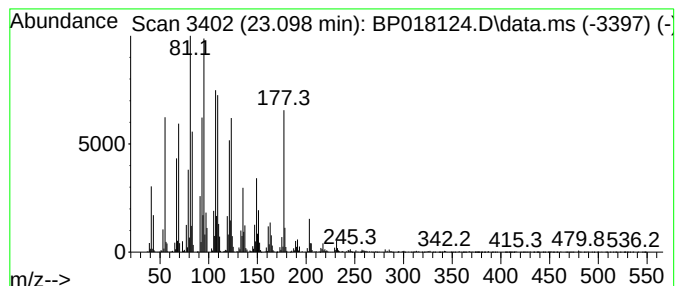
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 2-Methyl-3-(3-methyl-but-2-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.098	11.42 ng/ul	770152	Perylene-d12	23.927

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methyl-3-(3-methyl-but-2-enyl)...	222	C15H26O	1000144-10-2	94		
2	1-(2,2,3,5,6-Pentamethylcyclohex...	466	C34H58	1000373-94-7	91		
3	1,1,6-trimethyl-3-methylene-2-(3...	452	C33H56	1000373-94-5	72		
4	1,4-Dimethyl-7-(prop-1-en-2-yl)d...	222	C15H26O	021698-41-9	55		
5	(3S,6S)-6-Isopropyl-3-methyl-2-(...	220	C15H24O	021698-46-4	41		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018124.D
Acq On : 13 Nov 2023 02:47
Operator : MA/JU
Sample : 05158-02
Misc :
ALS Vial : 105 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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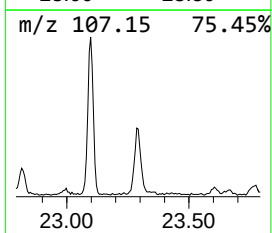
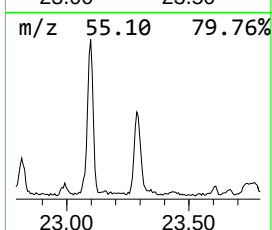
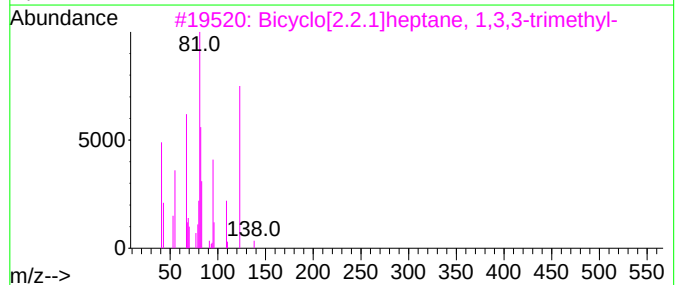
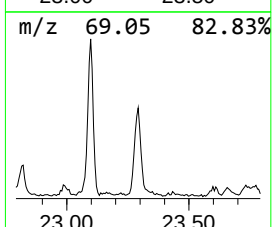
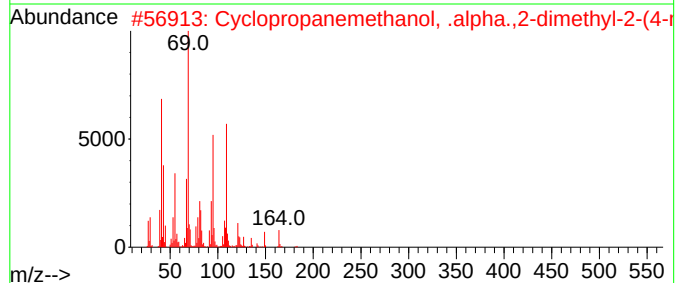
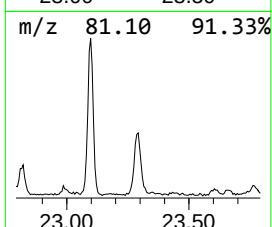
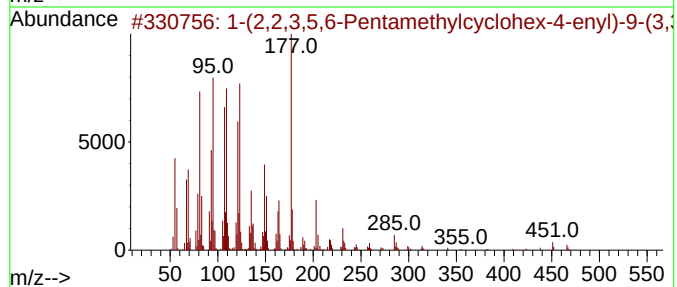
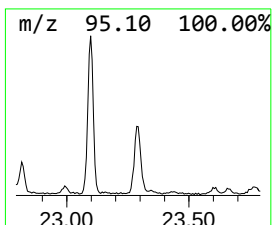
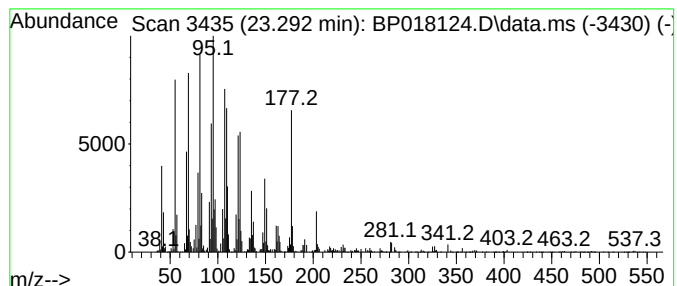
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 unknown-02 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.292	5.46 ng/ul	368329	Perylene-d12	23.927

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-(2,2,3,5,6-Pentamethylcyclohex...	466	C34H58	1000373-94-7	74
2			Cyclopropanemethanol, .alpha.,2-...	182	C12H22O	121959-70-4	22
3			Bicyclo[2.2.1]heptane, 1,3,3-tri...	138	C10H18	006248-88-0	18
4			Bicyclo[2.2.2]oct-2-ene, 1,2,3,6...	164	C12H20	062376-14-1	15
5			Vinyltris(cyclopropyl)silane	178	C11H18Si	1000109-97-3	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018124.D
Acq On : 13 Nov 2023 02:47
Operator : MA/JU
Sample : 05158-02
Misc :
ALS Vial : 105 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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1,6,10,14-Hexad...	18.633	2.9	ng/ul	188055	4	17.298	1296560	20.0
n-Nonadecanol-1	21.086	4.3	ng/ul	260158	5	21.427	1205230	20.0
unknown-01	22.163	2.3	ng/ul	138470	5	21.427	1205230	20.0
1-(2,2,3,5,6-Pe...	22.292	14.7	ng/ul	886250	5	21.427	1205230	20.0
1,1,6-trimethyl...	22.568	5.9	ng/ul	356825	5	21.427	1205230	20.0
Carbonic acid, ...	22.815	2.1	ng/ul	140760	6	23.927	1349110	20.0
2-Methyl-3-(3-m...	23.098	11.4	ng/ul	770152	6	23.927	1349110	20.0
unknown-02	23.292	5.5	ng/ul	368329	6	23.927	1349110	20.0