

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
 Data File : BP018123.D
 Acq On : 13 Nov 2023 02:13
 Operator : MA/JU
 Sample : 05158-01
 Misc :
 ALS Vial : 104 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH362

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: BP018123.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	21	24	28	rBV2	12754	14198	0.42%	0.047%
2	3.669	96	99	107	rBV	66691	108845	3.20%	0.361%
3	3.963	145	149	153	rVB7	3704	6897	0.20%	0.023%
4	4.187	185	187	191	rBV5	3020	3921	0.12%	0.013%
5	4.569	246	252	253	rBV5	2312	4387	0.13%	0.015%
6	5.387	387	391	397	rBV2	12327	17283	0.51%	0.057%
7	6.769	625	626	635	rVB9	3811	6324	0.19%	0.021%
8	7.022	662	669	683	rBV	47299	121323	3.57%	0.402%
9	7.193	692	698	711	rBV	209784	350874	10.32%	1.163%
10	7.369	721	728	741	rBV	222038	407193	11.97%	1.350%
11	7.840	802	808	824	rVB	185795	323727	9.52%	1.073%
12	8.040	837	842	848	rVB10	1737	4028	0.12%	0.013%
13	8.575	925	933	948	rBV	150240	326005	9.59%	1.081%
14	9.045	1005	1013	1028	rBV	265448	506835	14.90%	1.680%
15	9.216	1038	1042	1052	rVB2	31558	50590	1.49%	0.168%
16	9.763	1127	1135	1148	rBV	227586	427185	12.56%	1.416%
17	9.940	1161	1165	1170	rVB8	2907	4639	0.14%	0.015%
18	10.045	1176	1183	1186	rBV8	2028	3754	0.11%	0.012%
19	10.292	1216	1225	1250	rBV	258885	598802	17.61%	1.985%
20	10.657	1279	1287	1298	rBV	243639	428121	12.59%	1.419%
21	10.845	1312	1319	1339	rBV	225510	512202	15.06%	1.698%
22	11.698	1458	1464	1465	rBV5	2270	3731	0.11%	0.012%
23	12.098	1528	1532	1536	rBV6	2474	4173	0.12%	0.014%
24	12.269	1556	1561	1567	rBV8	4685	7992	0.24%	0.026%
25	12.328	1567	1571	1585	rBV2	26814	74899	2.20%	0.248%
26	13.357	1738	1746	1751	rBV4	11882	22646	0.67%	0.075%
27	13.722	1804	1808	1812	rVB7	2721	3493	0.10%	0.012%
28	13.939	1837	1845	1851	rBV	648301	917750	26.99%	3.043%
29	14.151	1877	1881	1884	rVB5	2043	3547	0.10%	0.012%
30	14.216	1885	1892	1903	rVV	673917	988603	29.07%	3.278%
31	14.298	1903	1906	1910	rVV6	4189	7494	0.22%	0.025%
32	14.351	1910	1915	1918	rVB7	3190	6019	0.18%	0.020%
33	14.516	1935	1943	1954	rBV	357431	523891	15.41%	1.737%
34	14.604	1955	1958	1962	rVB5	3231	5071	0.15%	0.017%
35	14.845	1992	1999	2007	rBV5	13633	45109	1.33%	0.150%
36	15.022	2027	2029	2034	rVB6	3094	3785	0.11%	0.013%

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37	15.251	2066	2068	2073	rBV6	2312	4426	0.13%	0.015%
38	15.298	2073	2076	2080	rVB6	4569	4822	0.14%	0.016%
39	15.522	2106	2114	2127	rBV	857673	1304533	38.36%	4.325%
40	15.680	2135	2141	2152	rBV	257691	451891	13.29%	1.498%

41	15.886	2171	2176	2180	rVB7	2164	3859	0.11%	0.013%
42	16.045	2199	2203	2204	rBV4	2516	3924	0.12%	0.013%
43	16.069	2205	2207	2211	rVB5	2750	3483	0.10%	0.012%
44	16.257	2234	2239	2242	rBV7	3957	7705	0.23%	0.026%
45	16.392	2260	2262	2268	rVB7	3261	4362	0.13%	0.014%

46	16.657	2304	2307	2313	rBV8	5553	7268	0.21%	0.024%
47	16.786	2324	2329	2337	rBV2	22477	48301	1.42%	0.160%
48	16.845	2337	2339	2343	rBV5	4707	5832	0.17%	0.019%
49	16.974	2355	2361	2365	rVB9	2746	4980	0.15%	0.017%
50	17.098	2377	2382	2386	rBV7	5188	10741	0.32%	0.036%

51	17.257	2405	2409	2411	rBV2	31459	38911	1.14%	0.129%
52	17.298	2411	2416	2427	rVV	439014	649518	19.10%	2.154%
53	17.398	2427	2433	2450	rVV2	1018214	1545233	45.44%	5.123%
54	17.586	2461	2465	2469	rVB7	12372	14707	0.43%	0.049%
55	17.916	2519	2521	2526	rVB6	4970	5290	0.16%	0.018%

56	18.145	2555	2560	2568	rBV	128723	192356	5.66%	0.638%
57	18.233	2571	2575	2581	rVB	46401	65753	1.93%	0.218%
58	18.398	2601	2603	2606	rBV4	3122	3676	0.11%	0.012%
59	18.563	2627	2631	2637	rVB3	35784	43383	1.28%	0.144%
60	18.627	2638	2642	2647	rBV	114459	136282	4.01%	0.452%

61	18.704	2653	2655	2660	rVB6	8265	8978	0.26%	0.030%
62	18.769	2664	2666	2668	rBV3	3754	3642	0.11%	0.012%
63	19.221	2739	2743	2746	rBV3	16540	22215	0.65%	0.074%
64	19.298	2753	2756	2759	rVB2	16305	18794	0.55%	0.062%
65	19.386	2767	2771	2777	rBV3	53866	73049	2.15%	0.242%

66	19.651	2810	2816	2826	rBV	1530708	2003958	58.93%	6.644%
67	21.086	3056	3060	3069	rVB2	112241	169743	4.99%	0.563%
68	21.427	3113	3118	3125	rBV	575022	796428	23.42%	2.641%
69	21.898	3195	3198	3202	rVB3	95645	108256	3.18%	0.359%
70	21.945	3204	3206	3209	rVB3	48096	48239	1.42%	0.160%

71	22.062	3222	3226	3230	rBV3	136876	168147	4.94%	0.558%
72	22.162	3239	3243	3253	rVB	478656	638983	18.79%	2.119%
73	22.292	3259	3265	3270	rBV	2631449	3400615	100.00%	11.275%
74	22.568	3307	3312	3320	rVB	888199	1294142	38.06%	4.291%
75	22.815	3346	3354	3363	rVB	442513	792513	23.30%	2.628%

76	22.992	3379	3384	3390	rBV4	126079	220069	6.47%	0.730%
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Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

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Title : SVOA CALIBRATION

77	23.098	3395	3402	3411	rVB	1974969	3239855	95.27%	10.742%
78	23.286	3428	3434	3442	rBV	934424	1784396	52.47%	5.916%
79	23.604	3484	3488	3493	rVB5	90366	133709	3.93%	0.443%
80	23.662	3494	3498	3504	rVB7	81668	132197	3.89%	0.438%
81	23.756	3507	3514	3526	rVB2	1333373	2734682	80.42%	9.067%
82	23.921	3536	3542	3552	rVB	456421	961240	28.27%	3.187%

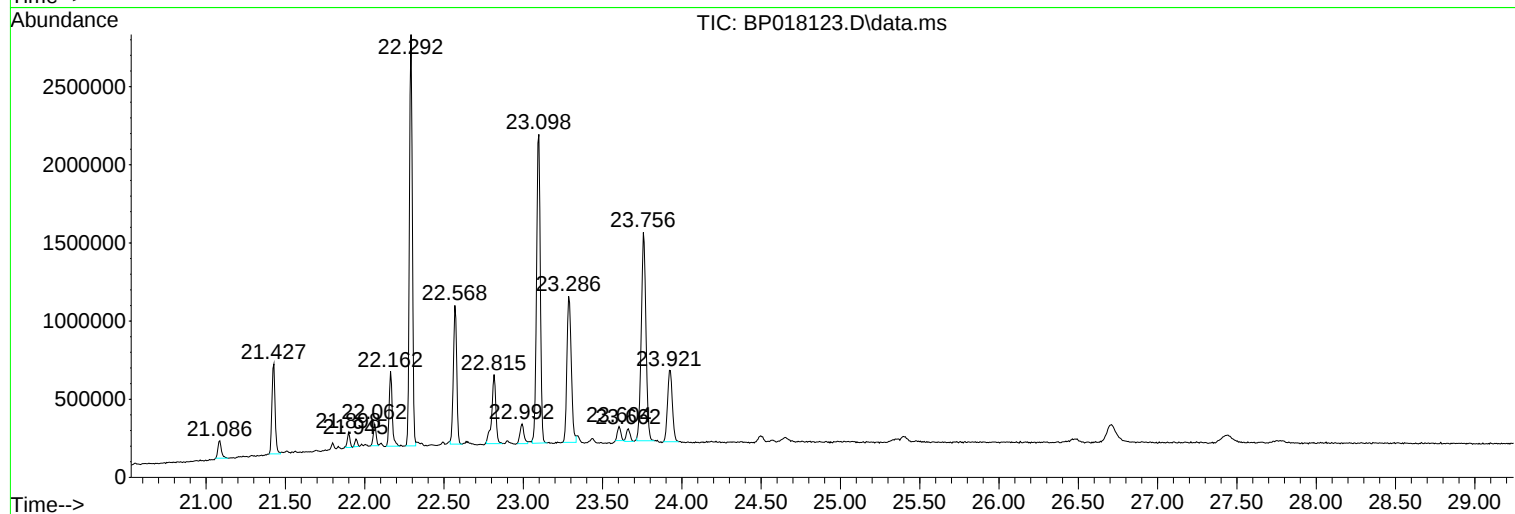
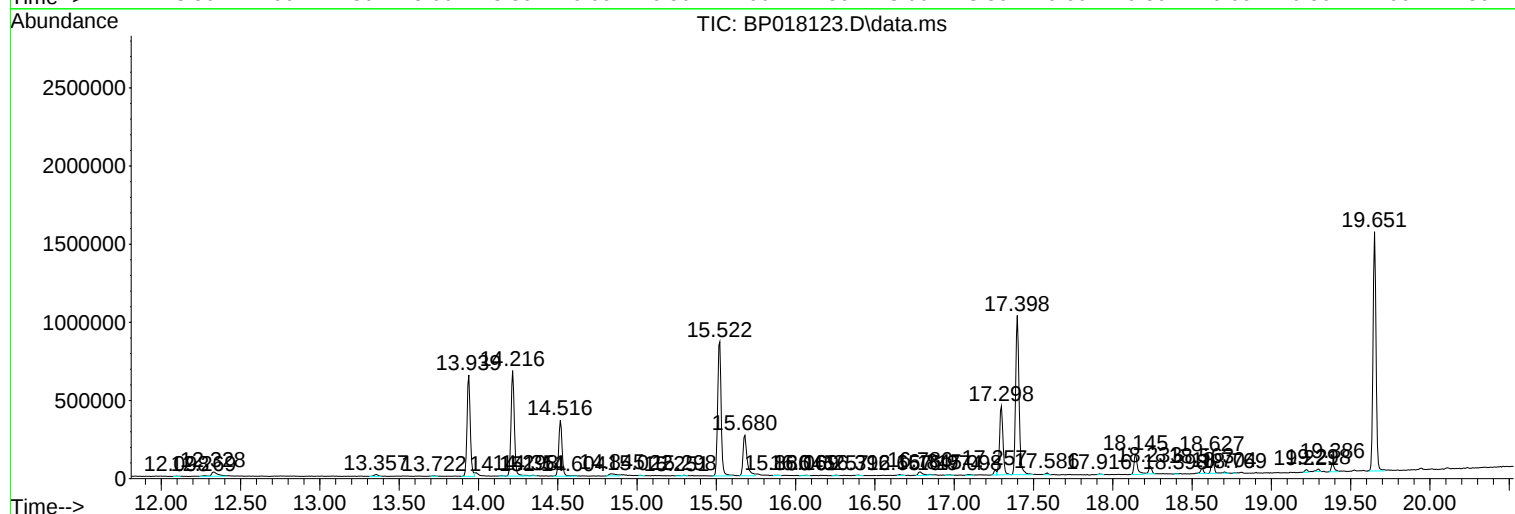
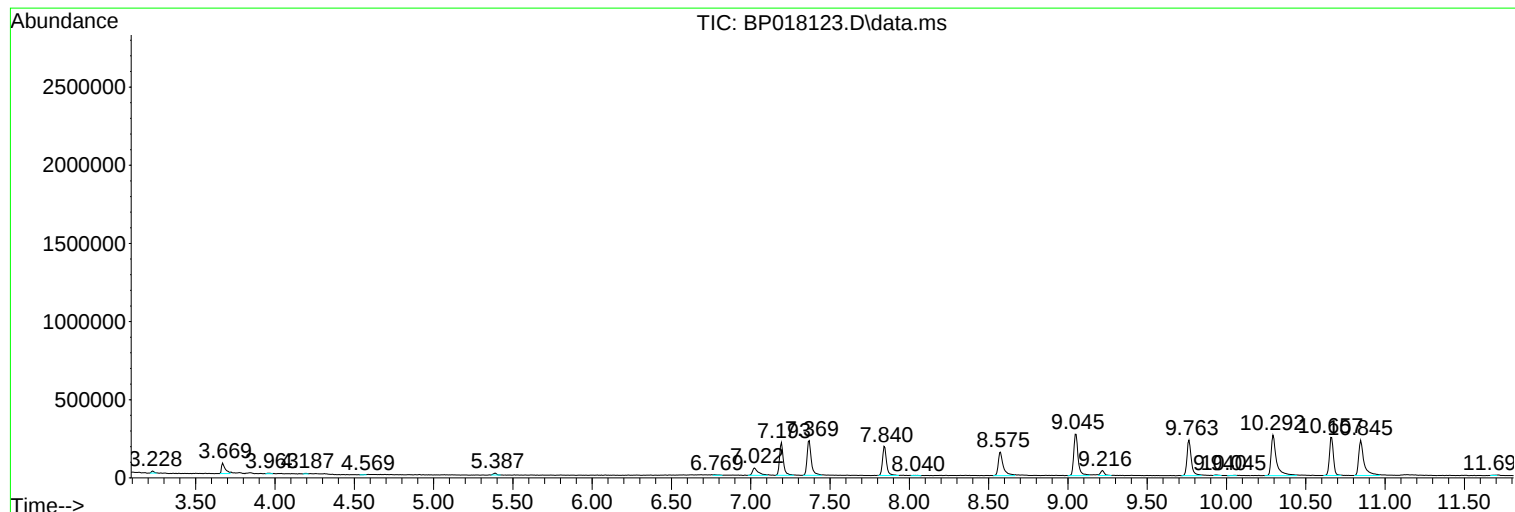
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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



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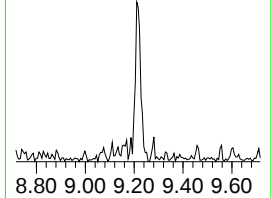
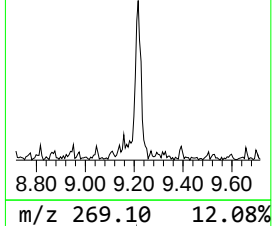
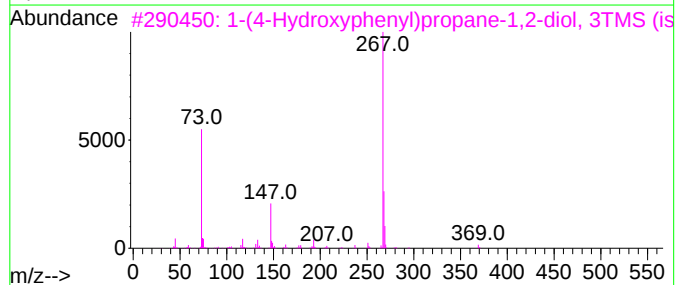
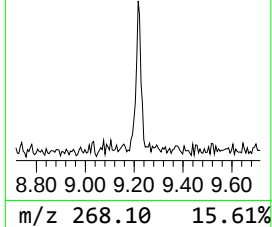
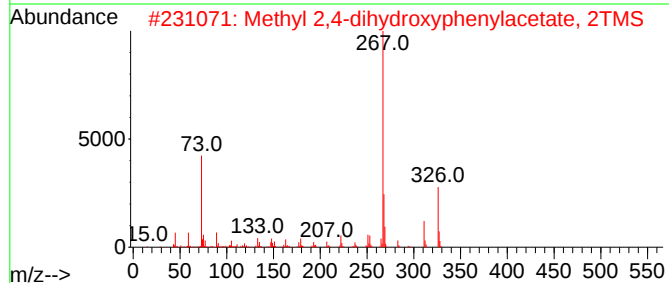
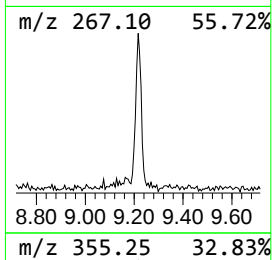
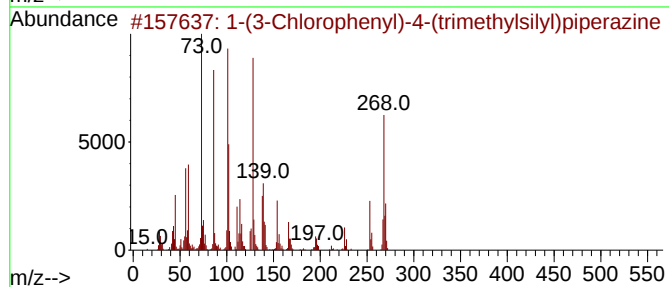
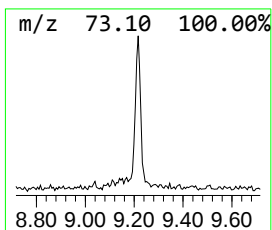
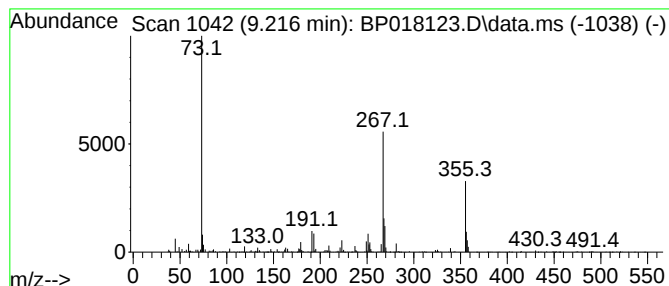
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 1-(3-Chlorophenyl)-4-(trime... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.216	3.13 ng/ul	50590	1,4-Dichlorobenzene-d4	7.840

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-(3-Chlorophenyl)-4-(trimethyls...	268	C13H21ClN2Si	1000408-33-9	53
2			Methyl 2,4-dihydroxyphenylacetat...	326	C15H26O4Si2	1010497-69-7	47
3			1-(4-Hydroxyphenyl)propane-1,2-d...	384	C18H36O3Si3	1000495-85-7	47
4			2-Amino-4-methylphenol, 2TMS der...	267	C13H25NOSi2	1000373-28-1	47
5			9-Acridinol, trimethylsilyl ether	267	C16H17NOSi	1000463-65-5	43



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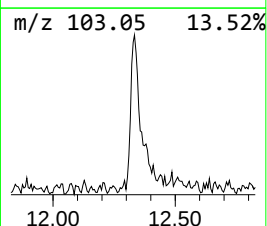
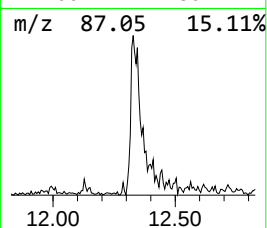
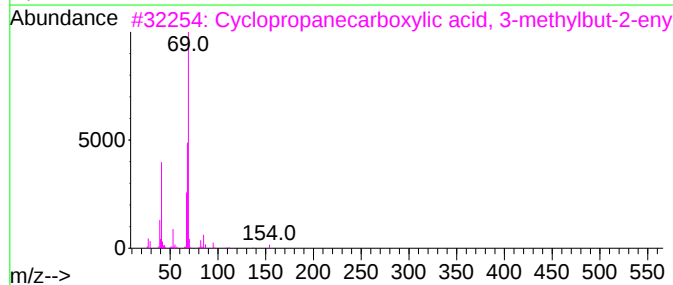
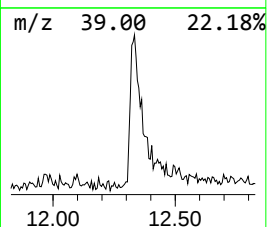
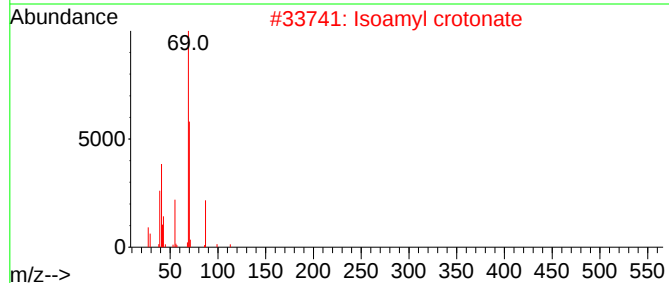
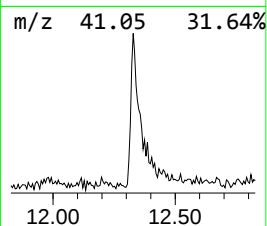
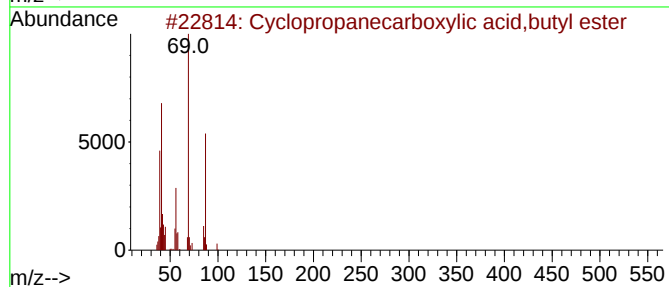
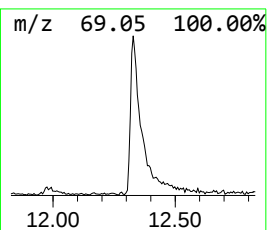
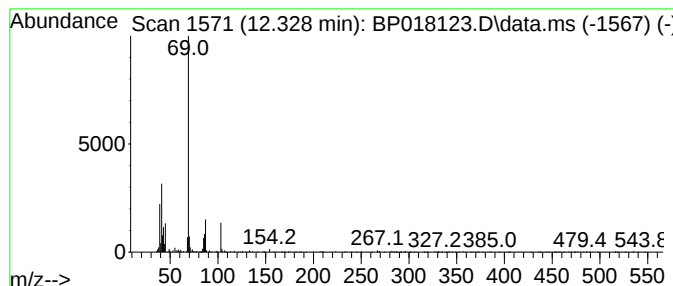
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Cyclopropanecarboxylic acid... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.328	3.50 ng/ul	74899	Naphthalene-d8	10.663

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropanecarboxylic acid, buty...	142	C8H14O2	054947-39-6	59
2			Isoamyl crotonate	156	C9H16O2	1010429-17-3	53
3			Cyclopropanecarboxylic acid, 3-m...	154	C9H14O2	1000299-37-4	50
4			2-Hydroxyethyl methacrylate	130	C6H10O3	000868-77-9	45
5			2-Butenoic acid, 2-propenyl ester	126	C7H10O2	020474-93-5	42



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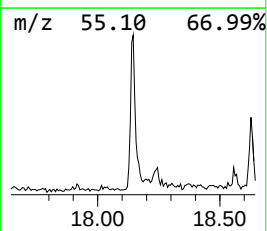
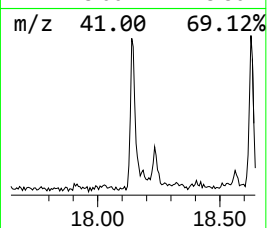
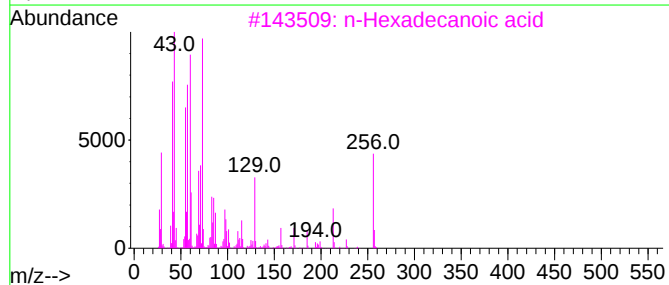
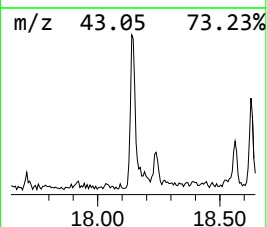
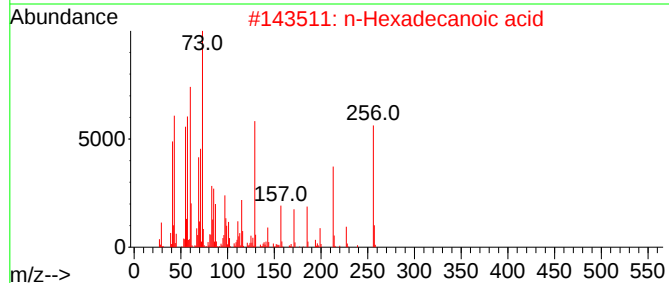
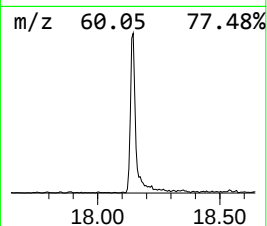
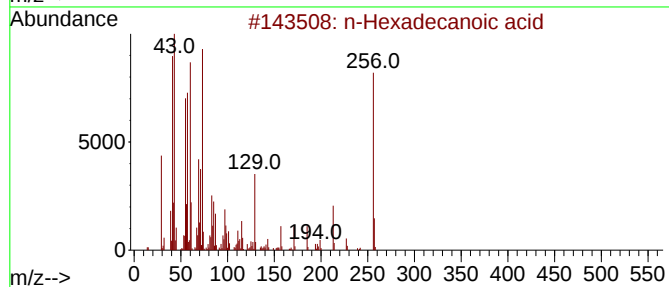
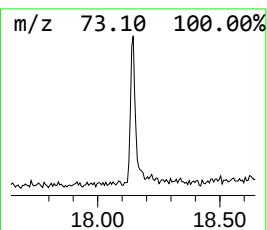
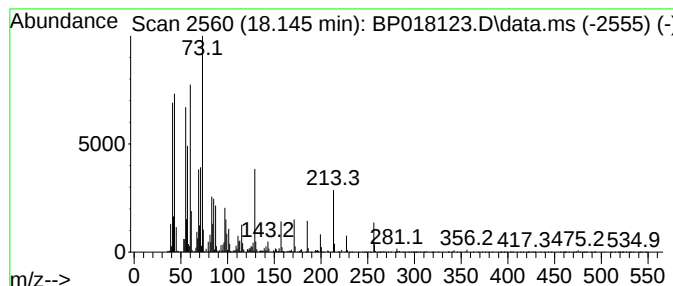
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.145	5.92 ng/ul	192356	Phenanthrene-d10	17.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	91



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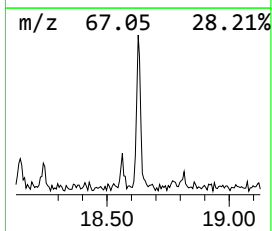
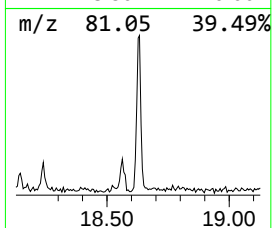
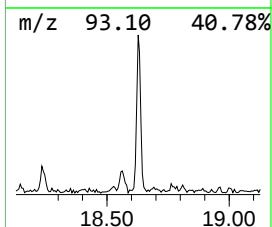
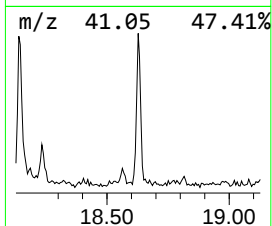
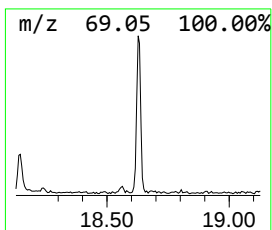
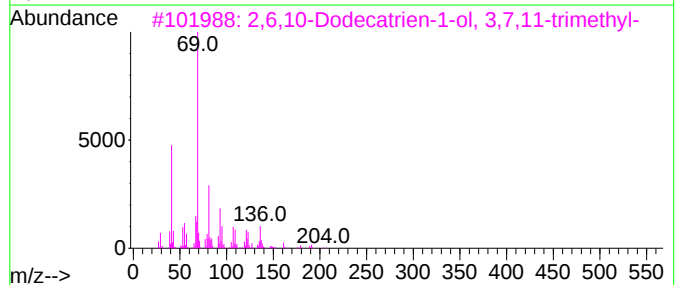
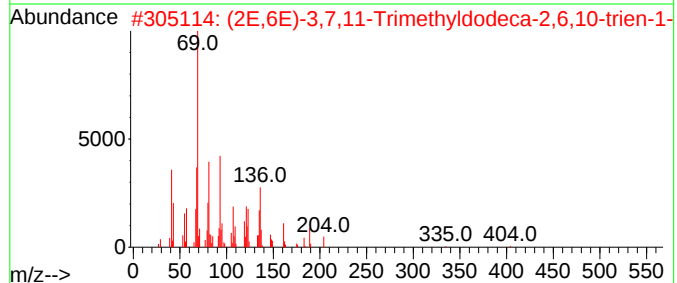
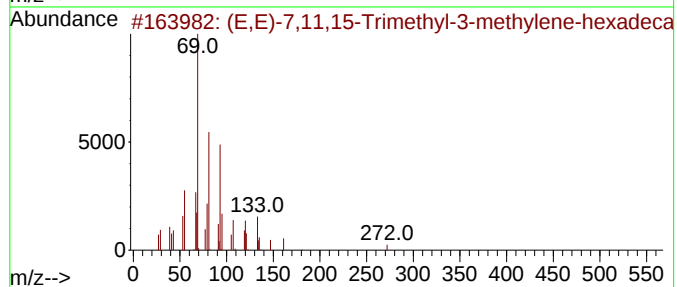
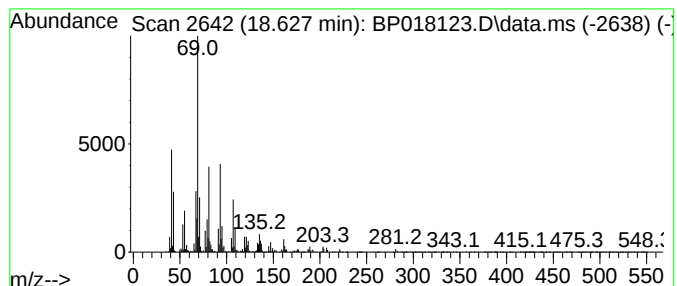
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (E,E)-7,11,15-Trimethyl-3-m... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.627	4.20 ng/ul	136282	Phenanthrene-d10	17.298

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E,E)-7,11,15-Trimethyl-3-methyl...	272	C20H32	070901-63-2	76		
2	(2E,6E)-3,7,11-Trimethyldodeca-2...	404	C27H48O2	078368-58-8	62		
3	2,6,10-Dodecatrien-1-ol, 3,7,11-...	222	C15H26O	004602-84-0	49		
4	2,6,10-Dodecatrien-1-ol, 3,7,11-...	222	C15H26O	003790-71-4	47		
5	Geranyl bromide	216	C10H17Br	006138-90-5	43		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018123.D
Acq On : 13 Nov 2023 02:13
Operator : MA/JU
Sample : 05158-01
Misc :
ALS Vial : 104 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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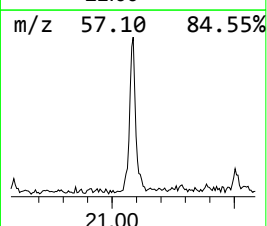
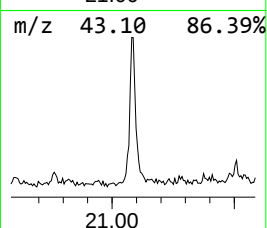
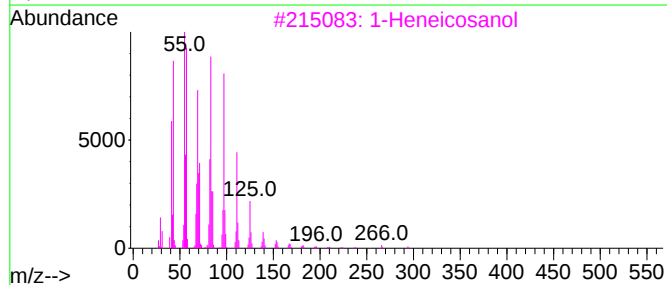
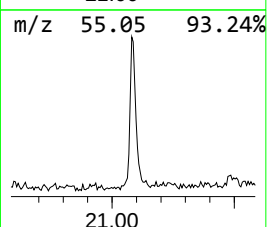
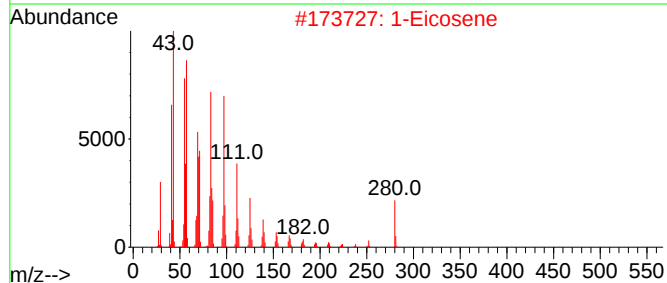
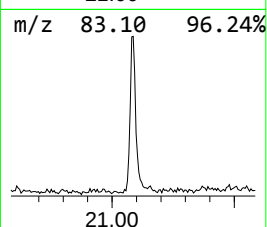
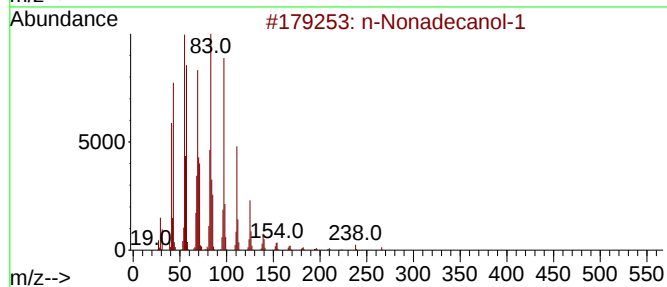
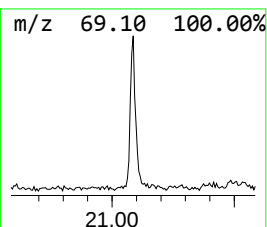
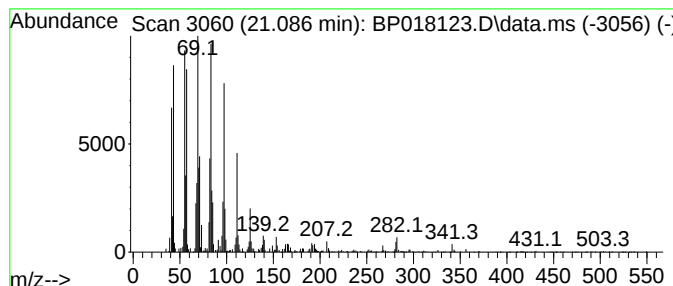
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 n-Nonadecanol-1 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.086	4.26 ng/ul	169743	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-Eicosene	280	C20H40	003452-07-1	91
3			1-Heneicosanol	312	C21H44O	015594-90-8	81
4			1-Octadecene	252	C18H36	000112-88-9	78
5			1-Eicosanol	298	C20H42O	000629-96-9	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018123.D
Acq On : 13 Nov 2023 02:13
Operator : MA/JU
Sample : 05158-01
Misc :
ALS Vial : 104 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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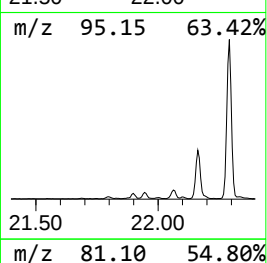
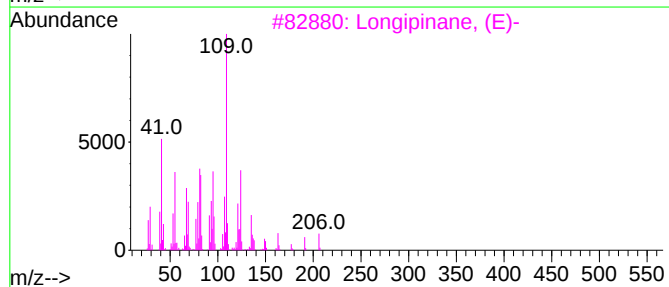
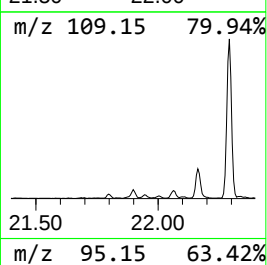
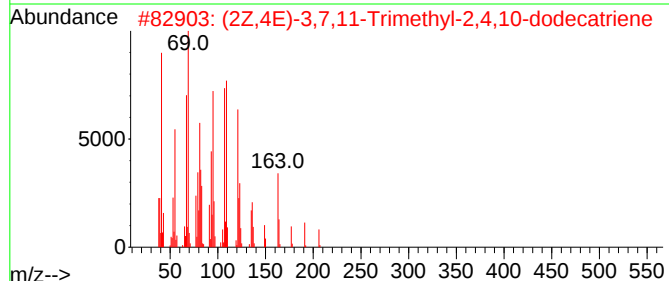
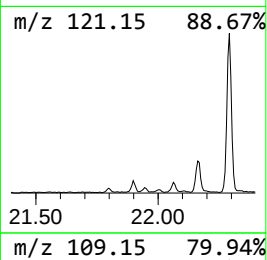
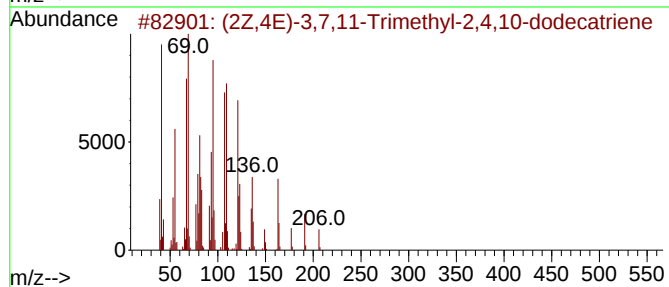
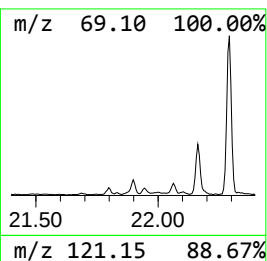
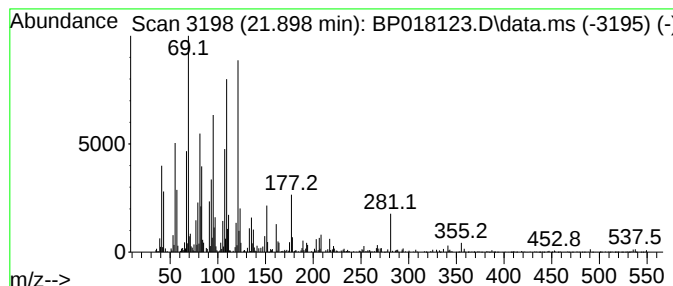
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-01 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.898	2.72 ng/ul	108256	Chrysene-d12	21.427

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(2Z,4E)-3,7,11-Trimethyl-2,4,10-...	206	C15H26	172549-29-0	30		
2	(2Z,4E)-3,7,11-Trimethyl-2,4,10-...	206	C15H26	172549-29-0	25		
3	Longipinane, (E)-	206	C15H26	1000156-14-5	20		
4	.alpha.-Estradiol, bis(trifluoro...	464	C22H22F6O4	1000352-22-8	18		
5	1-Methylene-2b-hydroxymethyl-3,3...	222	C15H26O	1000144-10-6	18		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018123.D
Acq On : 13 Nov 2023 02:13
Operator : MA/JU
Sample : 05158-01
Misc :
ALS Vial : 104 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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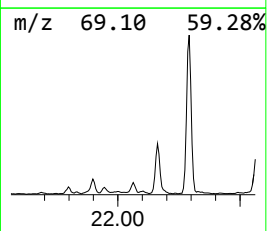
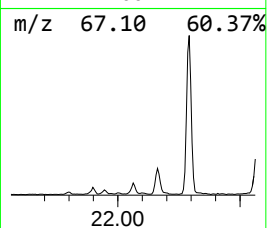
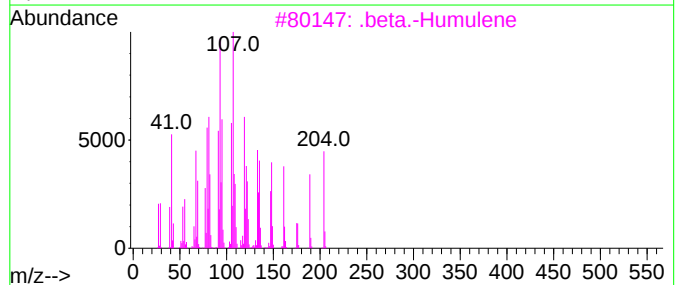
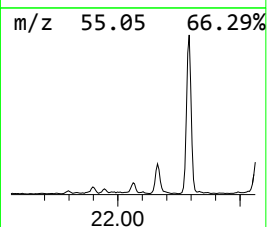
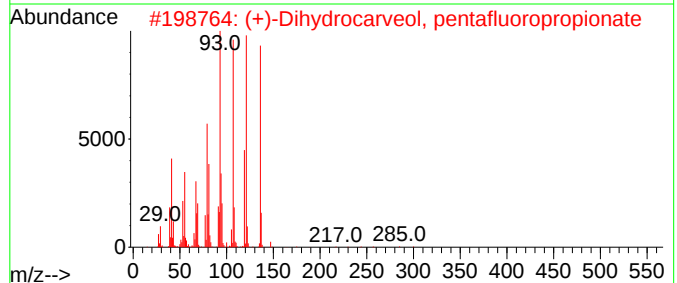
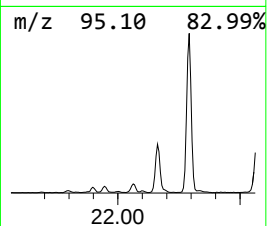
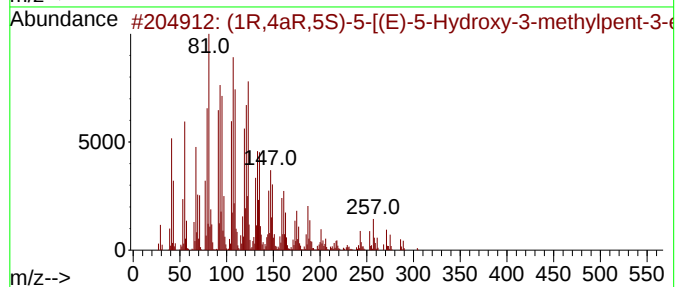
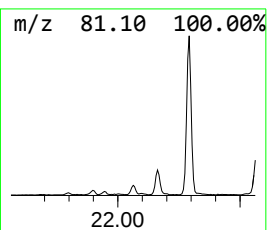
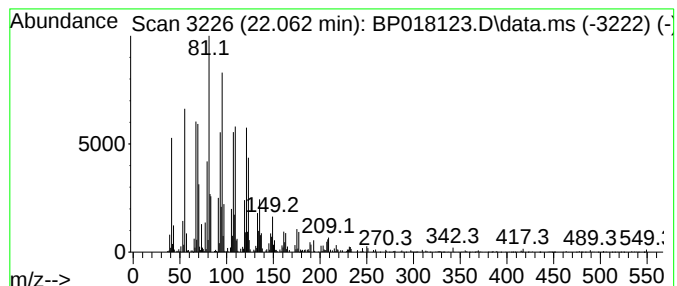
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (1R,4aR,5S)-5-[(E)-5-Hydrox... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.063	4.22 ng/ul	168147	Chrysene-d12	21.427

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(1R,4aR,5S)-5-[(E)-5-Hydroxy-3-m...	304	C20H32O2	1214984-94-7	60		
2	(+)-Dihydrocarveol, pentafluorop...	300	C13H17F5O2	1010462-97-4	42		
3	.beta.-Humulene	204	C15H24	000116-04-1	35		
4	(3R,3aS,6S,7R)-3,6,8,8-Tetrameth...	222	C15H26O	019903-73-2	22		
5	2-Buten-1-ol, 2-ethyl-4-(2,2,3-t...	208	C14H24O	028219-61-6	18		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018123.D
Acq On : 13 Nov 2023 02:13
Operator : MA/JU
Sample : 05158-01
Misc :
ALS Vial : 104 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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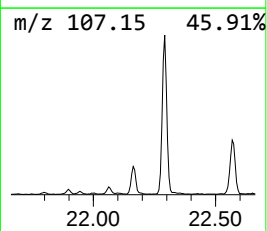
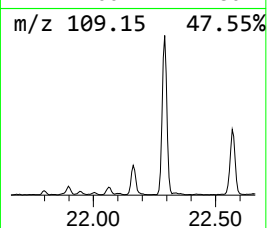
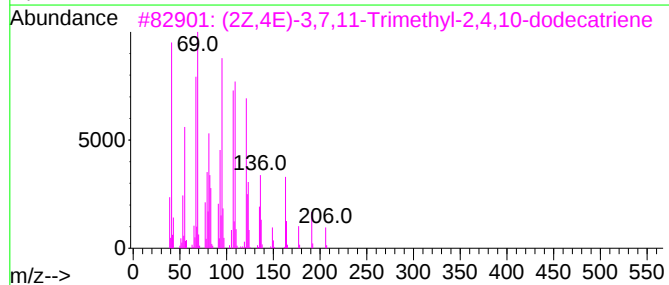
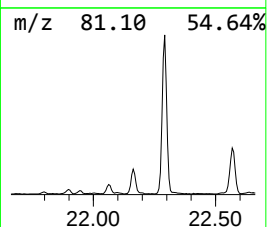
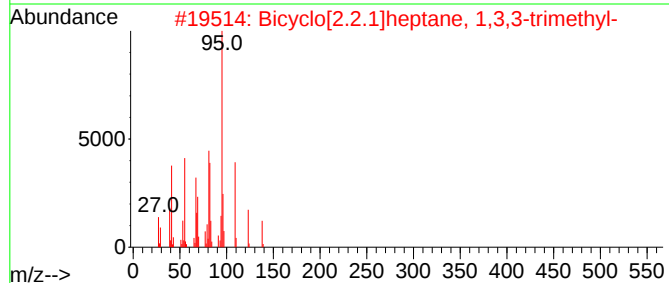
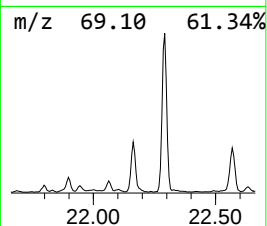
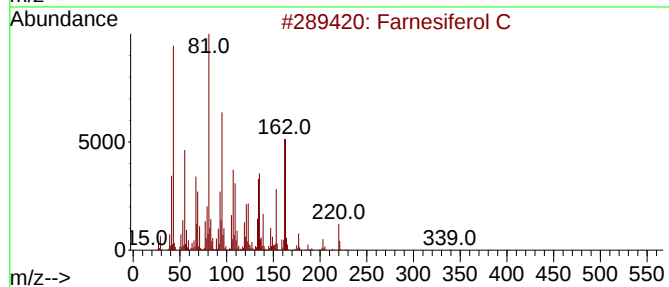
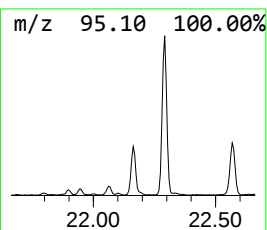
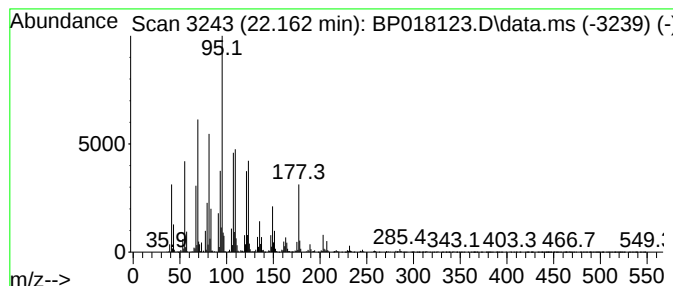
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Farnesiferol C Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.163	16.05 ng/ul	638983	Chrysene-d12	21.427

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Farnesiferol C	382	C24H30O4	000512-17-4	50
2			Bicyclo[2.2.1]heptane, 1,3,3-tri...	138	C10H18	006248-88-0	38
3			(2Z,4E)-3,7,11-Trimethyl-2,4,10-...	206	C15H26	172549-29-0	38
4			N-(2-Furylcarbonyl)tyrosine	275	C14H13NO5	1000260-99-5	35
5			N-(1,1-Dimethylpropynyl)-2-furan...	177	C10H11NO2	039108-97-9	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018123.D
Acq On : 13 Nov 2023 02:13
Operator : MA/JU
Sample : 05158-01
Misc :
ALS Vial : 104 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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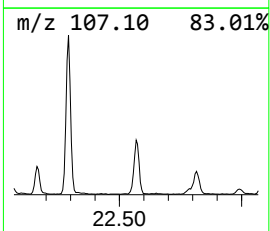
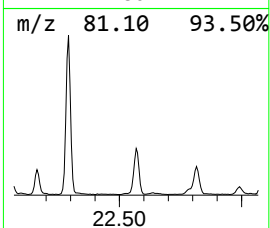
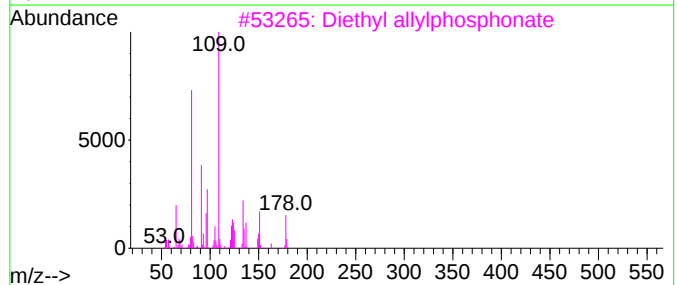
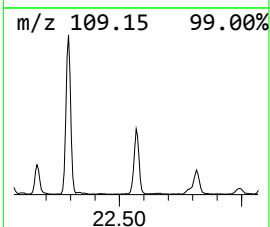
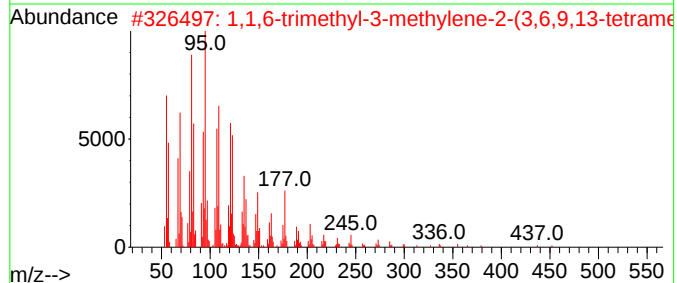
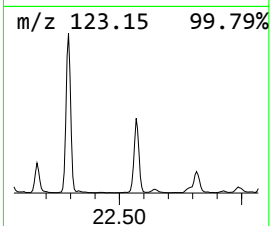
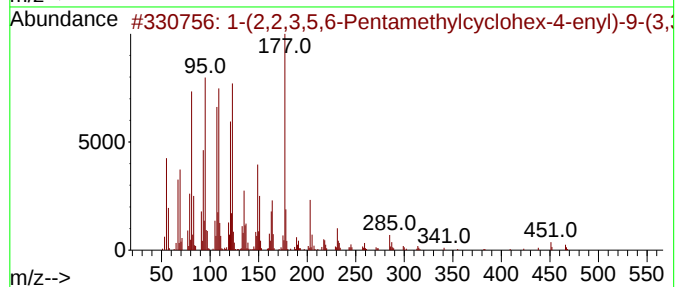
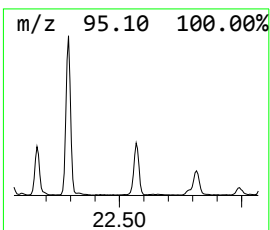
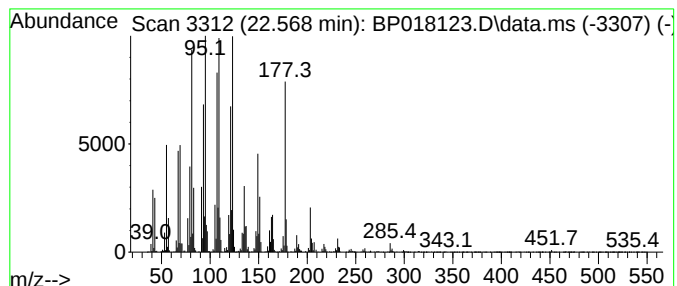
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-02 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.568	32.50 ng/ul	1294140	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	87		
2	1,1,6-trimethyl-3-methylene-2-(3...	452	C33H56	1000373-94-5	45		
3	Diethyl allylphosphonate	178	C7H15O3P	001067-87-4	18		
4	Camphorsulfonic acid	232	C10H16O4S	003144-16-9	15		
5	6-Methyl-6,7-dihydro-9H-5-oxa-9-...	177	C10H11NO2	1000210-20-2	14		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018123.D
Acq On : 13 Nov 2023 02:13
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Sample : 05158-01
Misc :
ALS Vial : 104 Sample Multiplier: 1

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ClientSampleId :
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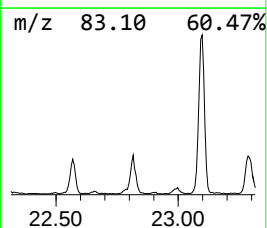
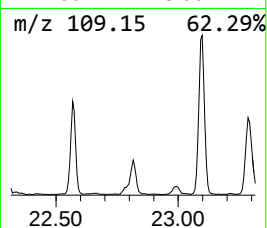
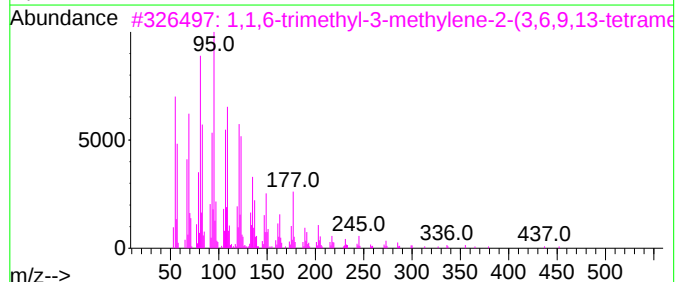
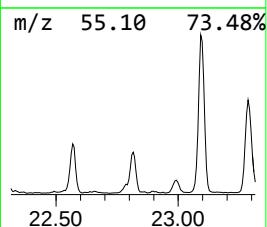
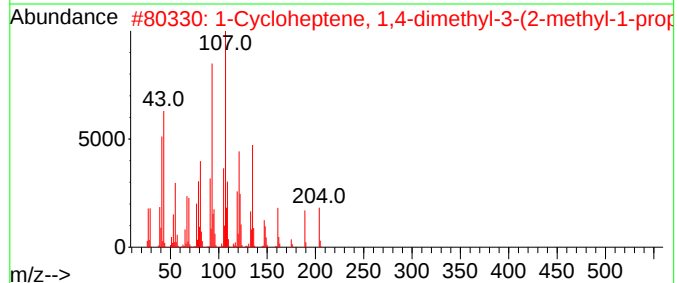
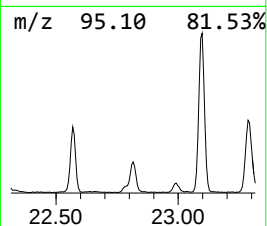
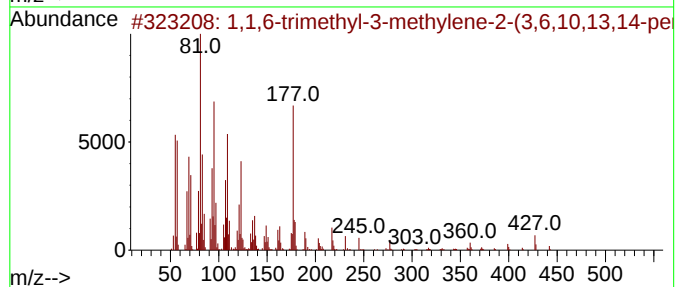
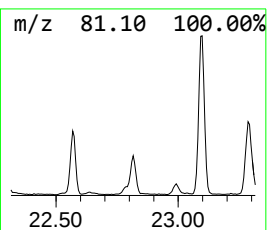
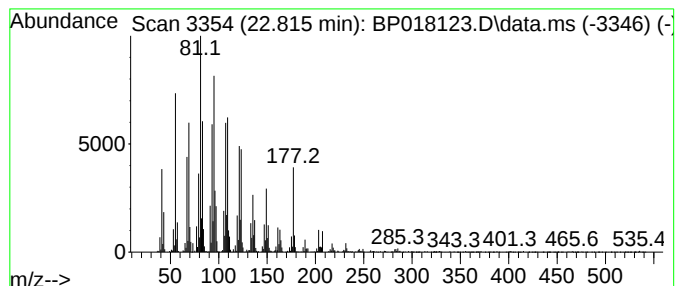
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 1,1,6-trimethyl-3-methylene... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.815	16.49 ng/ul	792513	Perylene-d12	23.921

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1,6-trimethyl-3-methylene-2-(3...	442	C32H58	1000373-94-0	64		
2	1-Cycloheptene, 1,4-dimethyl-3-(...	204	C15H24	1000159-38-6	56		
3	1,1,6-trimethyl-3-methylene-2-(3...	452	C33H56	1000373-94-5	46		
4	Vinyltris(cyclopropyl)silane	178	C11H18Si	1000109-97-3	27		
5	3-Octyne, 2,2,7-trimethyl-	152	C11H20	055402-13-6	27		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018123.D
Acq On : 13 Nov 2023 02:13
Operator : MA/JU
Sample : 05158-01
Misc :
ALS Vial : 104 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH362

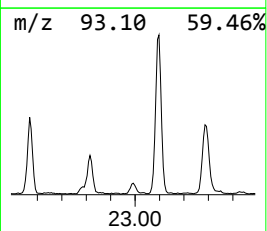
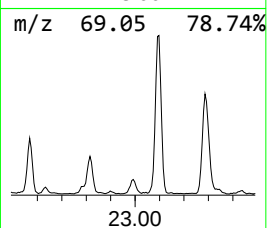
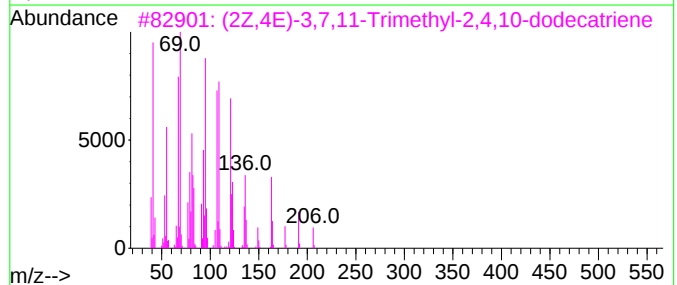
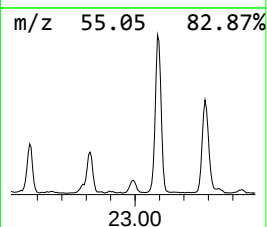
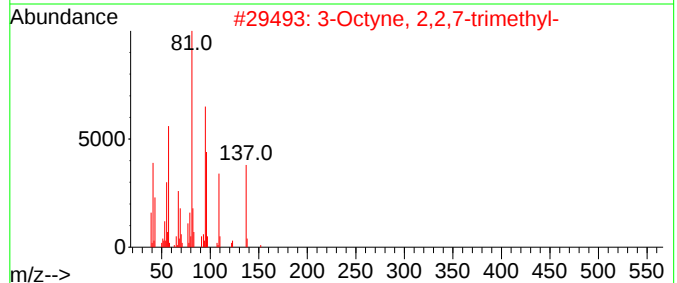
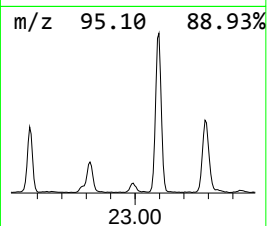
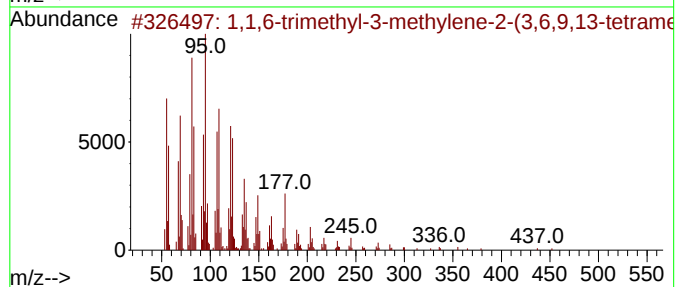
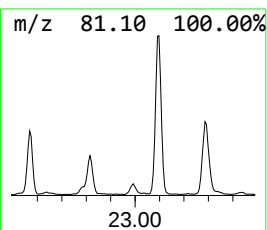
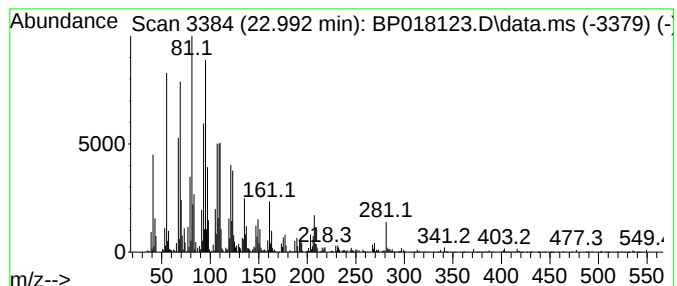
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 unknown-03 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.992	4.58 ng/ul	220069	Perylene-d12	23.921

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1,6-trimethyl-3-methylene-2-(3...	452	C33H56	1000373-94-5	46		
2	3-Octyne, 2,2,7-trimethyl-	152	C11H20	055402-13-6	43		
3	(2Z,4E)-3,7,11-Trimethyl-2,4,10-...	206	C15H26	172549-29-0	41		
4	22.alpha.-Hydroxy-3,4-secostict-...	458	C30H50O3	069773-95-1	38		
5	5-(1,2,4a,5-Tetramethyl-7-oxo-3,...	318	C20H30O3	1217871-71-0	25		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018123.D
Acq On : 13 Nov 2023 02:13
Operator : MA/JU
Sample : 05158-01
Misc :
ALS Vial : 104 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH362

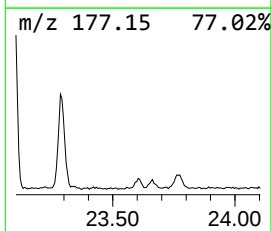
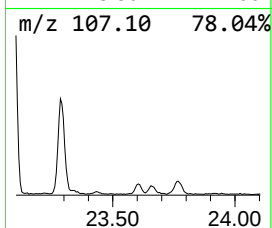
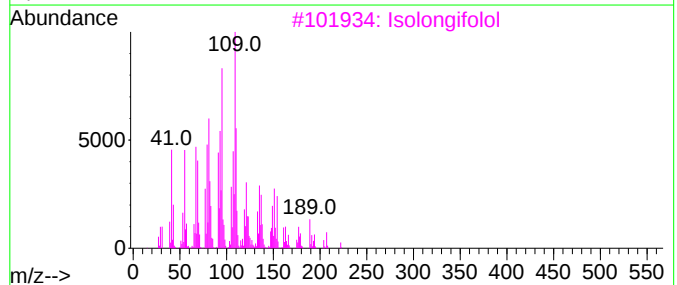
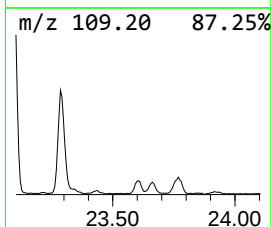
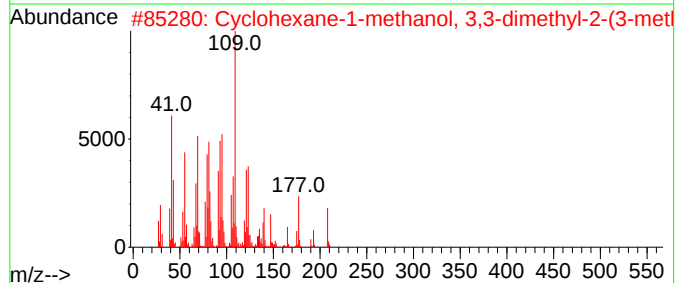
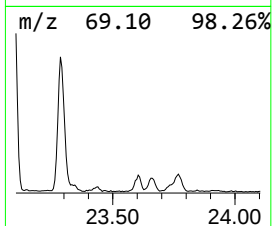
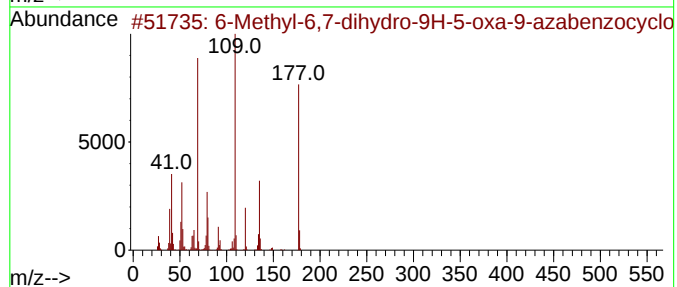
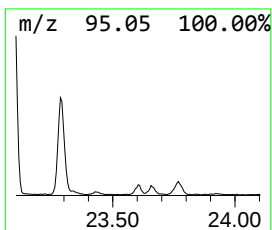
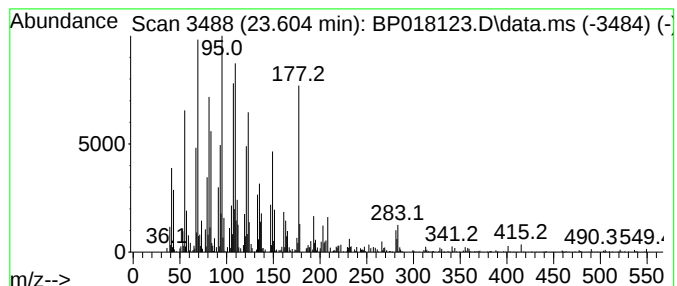
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 unknown-06 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.604	2.78 ng/ul	133709	Perylene-d12	23.921

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Methyl-6,7-dihydro-9H-5-oxa-9-...	177	C10H11NO2	1000210-20-2	43		
2	Cyclohexane-1-methanol, 3,3-dime...	208	C14H24O	1000196-01-5	38		
3	Isolongifolol	222	C15H26O	001139-17-9	35		
4	3-Chloro-2-trimethylsiloxypropene	164	C6H13ClOSi	1000427-07-8	20		
5	Cyclohexene, 1,5,5-trimethyl-6-a...	180	C12H20O	211563-96-1	14		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018123.D
Acq On : 13 Nov 2023 02:13
Operator : MA/JU
Sample : 05158-01
Misc :
ALS Vial : 104 Sample Multiplier: 1

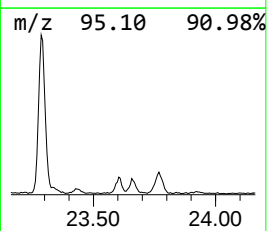
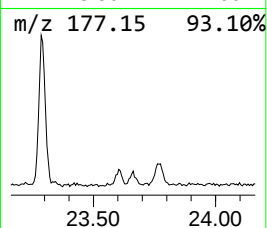
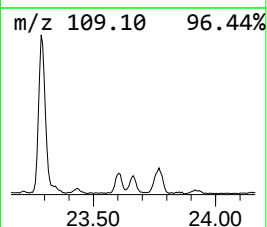
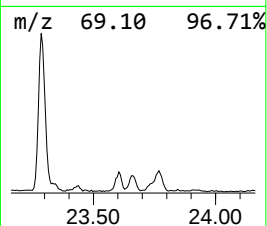
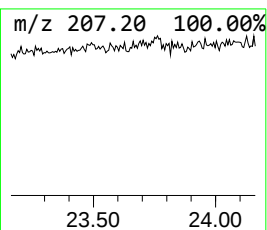
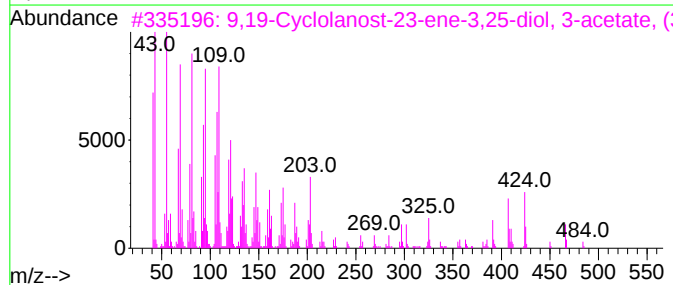
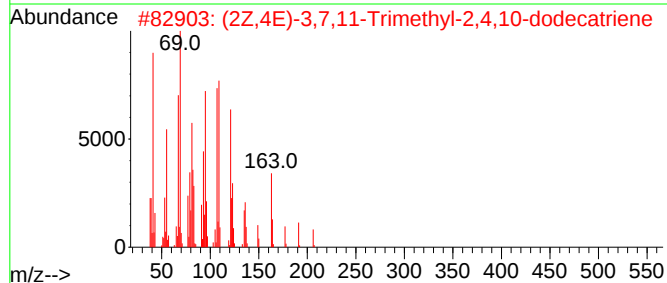
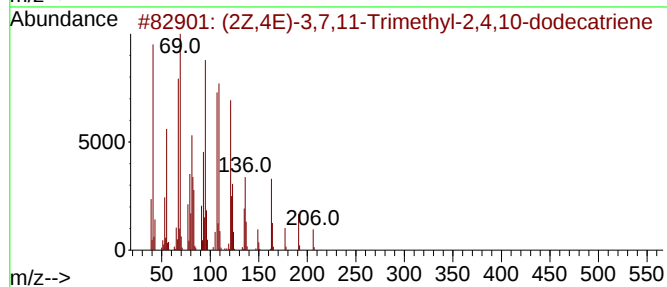
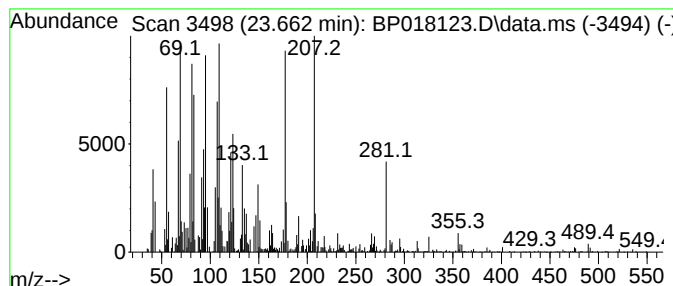
Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 16 unknown-07 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.662	2.75 ng/ul	132197	Perylene-d12	23.921	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(2Z,4E)-3,7,11-Trimethyl-2,4,10-...	206	C15H26	172549-29-0	46
2	(2Z,4E)-3,7,11-Trimethyl-2,4,10-...	206	C15H26	172549-29-0	38
3	9,19-Cyclolanost-23-ene-3,25-dio...	484	C32H52O3	054482-56-3	37
4	2H-3,9a-Methano-1-benzoxepin, oc...	222	C15H26O	005956-09-2	35
5	1-Formyl-2,2-dimethyl-3-trans-(3...	220	C15H24O	1000144-09-7	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
Data File : BP018123.D
Acq On : 13 Nov 2023 02:13
Operator : MA/JU
Sample : 05158-01
Misc :
ALS Vial : 104 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH362

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
1-(3-Chlorophen...	9.216	3.1	ng/ul	50590	1	7.840	323727	20.0
Cyclopropanecar...	12.328	3.5	ng/ul	74899	2	10.663	428121	20.0
n-Hexadecanoic ...	18.145	5.9	ng/ul	192356	4	17.298	649518	20.0
(E,E)-7,11,15-T...	18.627	4.2	ng/ul	136282	4	17.298	649518	20.0
n-Nonadecanol-1	21.086	4.3	ng/ul	169743	5	21.427	796428	20.0
unknown-01	21.898	2.7	ng/ul	108256	5	21.427	796428	20.0
(1R,4aR,5S)-5-[...	22.063	4.2	ng/ul	168147	5	21.427	796428	20.0
Farnesiferol C	22.163	16.1	ng/ul	638983	5	21.427	796428	20.0
1-(2,2,3,5,6-Pe...	22.292	85.4	ng/ul	3400620	5	21.427	796428	20.0
unknown-02	22.568	32.5	ng/ul	1294140	5	21.427	796428	20.0
1,1,6-trimethyl...	22.815	16.5	ng/ul	792513	6	23.921	961240	20.0
unknown-03	22.992	4.6	ng/ul	220069	6	23.921	961240	20.0
unknown-04	23.098	67.4	ng/ul	3239860	6	23.921	961240	20.0
unknown-05	23.286	37.1	ng/ul	1784400	6	23.921	961240	20.0
unknown-06	23.604	2.8	ng/ul	133709	6	23.921	961240	20.0
unknown-07	23.662	2.8	ng/ul	132197	6	23.921	961240	20.0