

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
 Data File : BM033944.D
 Acq On : 01 Feb 2022 17:56
 Operator : CG/JU
 Sample : N1273-21
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 COHL9

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
 Title : SVOA CALIBRATION

Signal : TIC: BM033944.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.325	11	15	27	rVB	695819	937718	18.92%	2.661%
2	3.578	55	58	66	rVB4	11621	19355	0.39%	0.055%
3	3.725	80	83	97	rBV	85483	139319	2.81%	0.395%
4	3.849	101	104	111	rVB2	9517	15457	0.31%	0.044%
5	4.060	136	140	144	rBV2	12734	19556	0.39%	0.056%
6	4.119	148	150	155	rVV3	8813	10534	0.21%	0.030%
7	4.184	157	161	165	rVV	22157	30844	0.62%	0.088%
8	4.231	165	169	174	rVB6	3261	5697	0.11%	0.016%
9	4.413	196	200	208	rVB6	7294	14090	0.28%	0.040%
10	4.607	226	233	245	rVV	3249421	4957184	100.00%	14.069%
11	4.701	245	249	254	rVB	8208	11429	0.23%	0.032%
12	5.348	354	359	366	rVB	25424	37133	0.75%	0.105%
13	6.925	620	627	630	rBV7	3047	5634	0.11%	0.016%
14	7.084	648	654	665	rVB2	103944	189941	3.83%	0.539%
15	7.248	676	682	692	rBV	413909	674811	13.61%	1.915%
16	7.454	710	717	727	rBV	390281	613596	12.38%	1.741%
17	7.719	757	762	769	rVB5	8292	11867	0.24%	0.034%
18	7.931	790	798	806	rBV	394852	631757	12.74%	1.793%
19	8.001	806	810	813	rBV2	6728	9311	0.19%	0.026%
20	8.636	911	918	929	rBV2	284756	457465	9.23%	1.298%
21	8.748	929	937	942	rVB6	4155	9411	0.19%	0.027%
22	9.089	988	995	1006	rVB	569486	905646	18.27%	2.570%
23	9.207	1011	1015	1022	rVB7	3579	6263	0.13%	0.018%
24	9.325	1032	1035	1040	rVB6	4111	4990	0.10%	0.014%
25	9.542	1066	1072	1080	rVB	89266	136480	2.75%	0.387%
26	9.819	1110	1119	1133	rBV	405623	685814	13.83%	1.946%
27	10.360	1203	1211	1223	rBV	559495	927344	18.71%	2.632%
28	10.519	1232	1238	1250	rBV	71060	124650	2.51%	0.354%
29	10.689	1262	1267	1269	rBV3	4759	7063	0.14%	0.020%
30	10.742	1269	1276	1284	rVB	550144	919380	18.55%	2.609%
31	10.813	1284	1288	1291	rBV3	7833	11760	0.24%	0.033%
32	10.872	1291	1298	1310	rBV	520513	893865	18.03%	2.537%
33	12.036	1490	1496	1502	rBV2	62034	94162	1.90%	0.267%
34	12.195	1520	1523	1529	rVB2	7995	13021	0.26%	0.037%
35	12.330	1541	1546	1551	rVB	11346	18799	0.38%	0.053%
36	12.760	1613	1619	1623	rBV7	3132	5064	0.10%	0.014%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
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37	13.089	1666	1675	1678	rBV8	3544	8924	0.18%	0.025%
38	13.177	1685	1690	1697	rVB7	6546	12391	0.25%	0.035%
39	13.413	1723	1730	1736	rBV2	134745	211192	4.26%	0.599%
40	13.466	1736	1739	1747	rVB2	9301	15293	0.31%	0.043%
41	13.889	1802	1811	1816	rBV4	17034	32022	0.65%	0.091%
42	13.977	1819	1826	1831	rBV	1297004	1802994	36.37%	5.117%
43	14.013	1831	1832	1838	rVV2	27827	26564	0.54%	0.075%
44	14.177	1852	1860	1868	rBV	223505	415219	8.38%	1.178%
45	14.260	1868	1874	1889	rVB	1415208	2070825	41.77%	5.877%
46	14.471	1906	1910	1914	rVB	9018	12113	0.24%	0.034%
47	14.566	1917	1926	1935	rVV	832523	1237678	24.97%	3.513%
48	14.760	1954	1959	1970	rVV2	48474	103522	2.09%	0.294%
49	14.948	1981	1991	1997	rBV6	23049	44764	0.90%	0.127%
50	15.195	2028	2033	2038	rBV4	7357	11532	0.23%	0.033%
51	15.307	2048	2052	2056	rVB	8302	10661	0.22%	0.030%
52	15.365	2057	2062	2066	rVV3	6670	11518	0.23%	0.033%
53	15.418	2068	2071	2077	rVB2	7544	9255	0.19%	0.026%
54	15.554	2087	2094	2106	rBV	1865238	2591785	52.28%	7.356%
55	15.665	2107	2113	2121	rVV	544082	730691	14.74%	2.074%
56	15.789	2130	2134	2138	rVB5	12693	21369	0.43%	0.061%
57	15.995	2166	2169	2174	rVB6	6637	9253	0.19%	0.026%
58	16.195	2197	2203	2206	rBV7	4465	6665	0.13%	0.019%
59	16.236	2206	2210	2213	rBV6	6759	8916	0.18%	0.025%
60	16.277	2213	2217	2219	rBV4	6652	8430	0.17%	0.024%
61	16.330	2224	2226	2231	rVB6	4409	6052	0.12%	0.017%
62	16.442	2241	2245	2251	rBV6	7540	13044	0.26%	0.037%
63	16.807	2304	2307	2311	rVB2	13211	16652	0.34%	0.047%
64	16.912	2321	2325	2329	rBV7	5163	9795	0.20%	0.028%
65	16.954	2329	2332	2335	rVB3	8554	10894	0.22%	0.031%
66	17.065	2348	2351	2354	rBV3	8633	11456	0.23%	0.033%
67	17.224	2375	2378	2381	rVB5	6228	6901	0.14%	0.020%
68	17.301	2385	2391	2401	rBV	974076	1366773	27.57%	3.879%
69	17.401	2401	2408	2422	rVV	1905214	2747657	55.43%	7.798%
70	17.806	2473	2477	2481	rBV6	4099	5941	0.12%	0.017%
71	17.877	2485	2489	2493	rVB2	11550	14596	0.29%	0.041%
72	17.977	2501	2506	2510	rBV	92110	112585	2.27%	0.320%
73	18.089	2518	2525	2528	rBV9	5451	12567	0.25%	0.036%
74	18.195	2538	2543	2548	rBV3	52567	73946	1.49%	0.210%
75	18.495	2592	2594	2600	rBV6	3085	5716	0.12%	0.016%
76	18.665	2618	2623	2626	rBV7	3420	7355	0.15%	0.021%

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77	18.818	2645	2649	2655	rBV4	13034	18949	0.38%	0.054%
78	19.142	2700	2704	2708	rVV	89114	100191	2.02%	0.284%
79	19.253	2719	2723	2726	rBV3	24732	35743	0.72%	0.101%
80	19.318	2731	2734	2738	rVV2	21030	28141	0.57%	0.080%
81	19.465	2756	2759	2763	rBV4	17428	22462	0.45%	0.064%
82	19.606	2780	2783	2788	rVB3	27605	38636	0.78%	0.110%
83	19.683	2790	2796	2803	rBV	2051893	2827056	57.03%	8.023%
84	21.177	3047	3050	3057	rBV5	46810	68688	1.39%	0.195%
85	21.465	3094	3099	3104	rBV	942512	1176410	23.73%	3.339%
86	23.700	3472	3479	3486	rBV2	1232150	2399336	48.40%	6.809%
87	23.853	3499	3505	3514	rVB2	579933	1128466	22.76%	3.203%

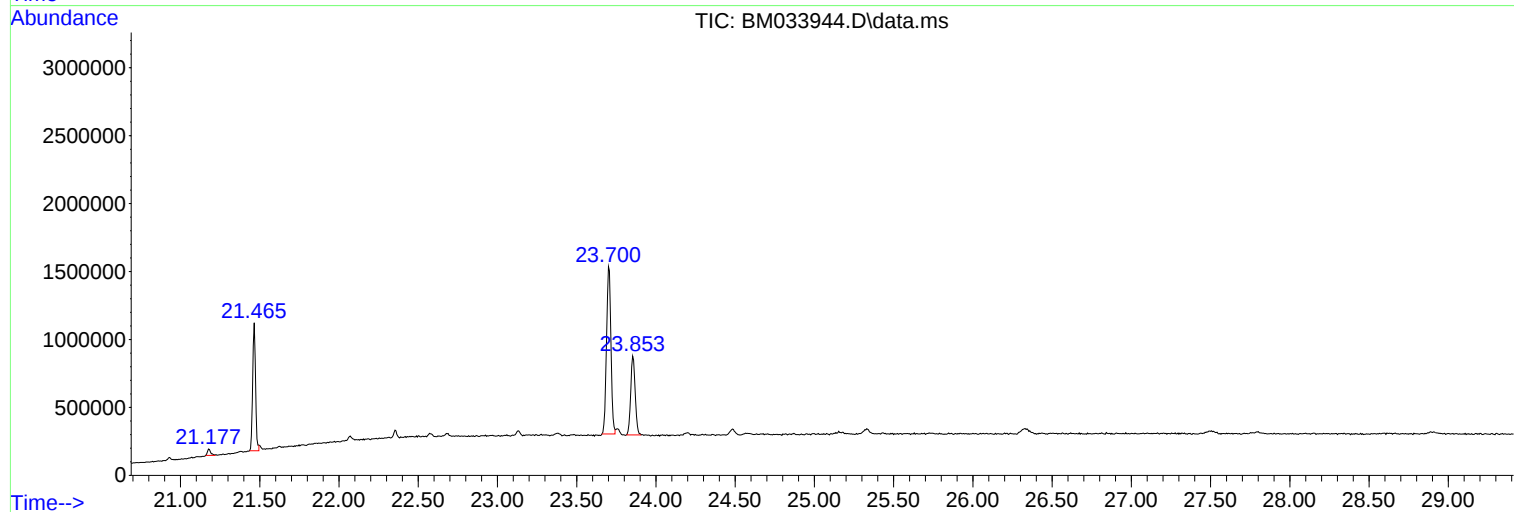
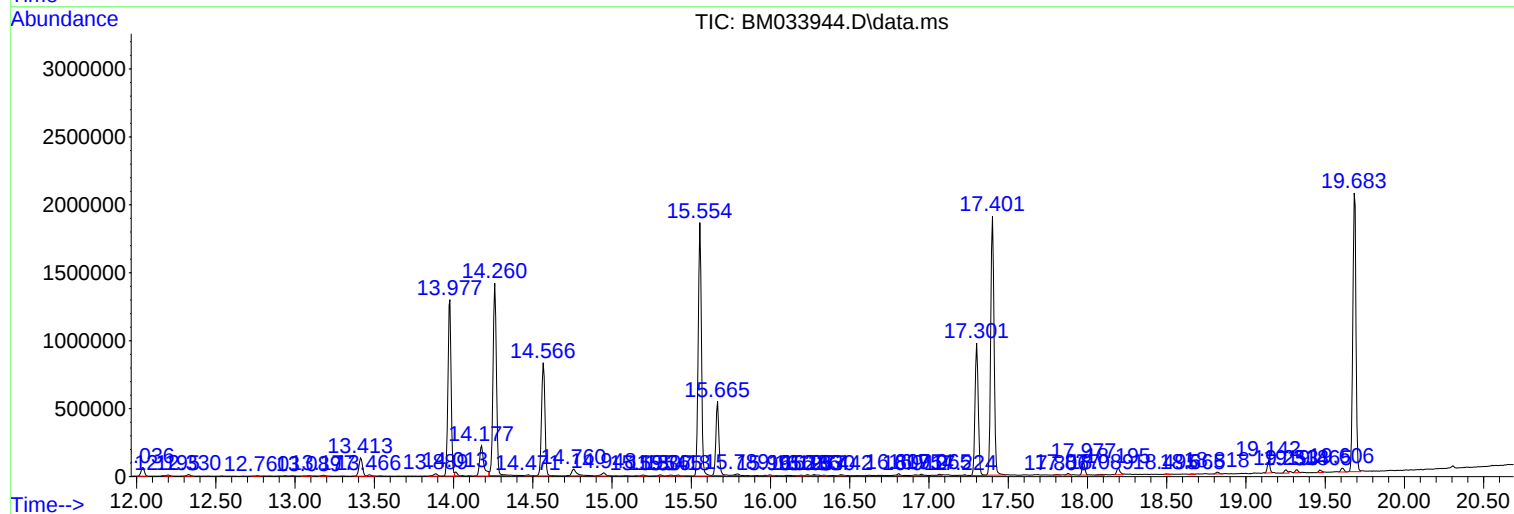
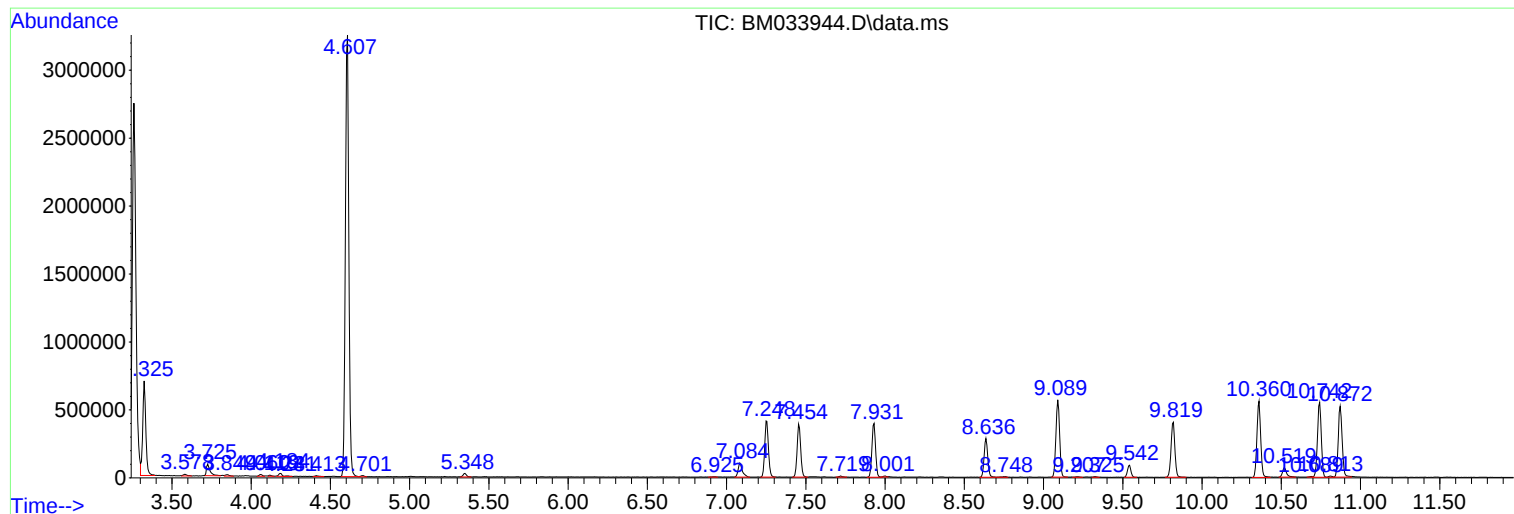
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

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



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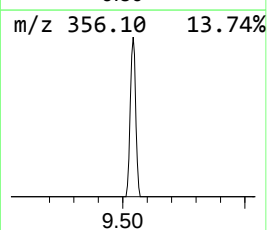
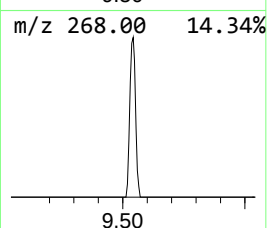
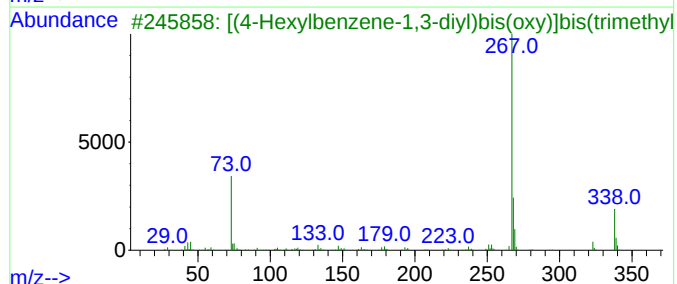
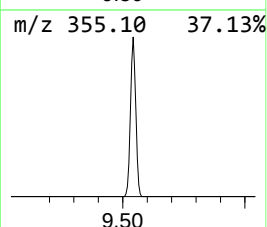
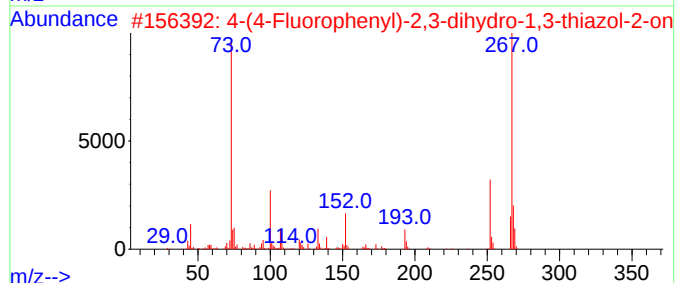
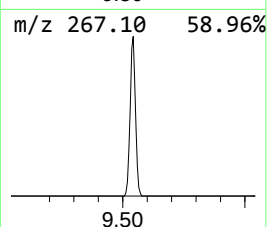
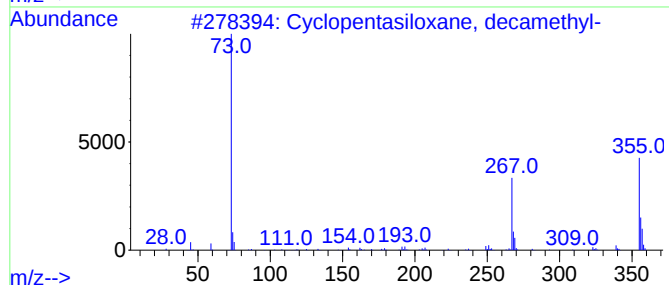
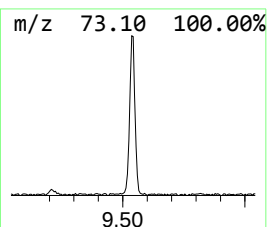
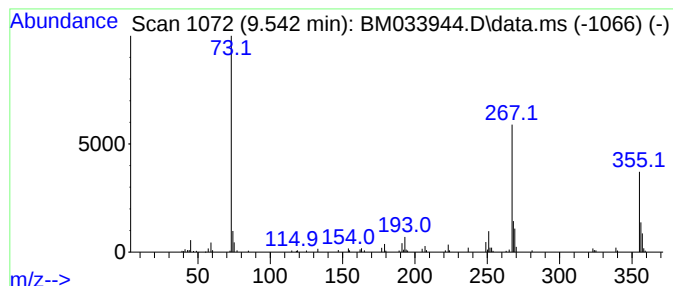
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Cyclopentasiloxane, decamet... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.542	2.97 ng/ul	136480	Naphthalene-d8	10.742

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentasiloxane, decamethyl-	370	C10H30O5Si5	000541-02-6	52
2			4-(4-Fluorophenyl)-2,3-dihydro-1...	267	C12H14FNOSSi	1010501-61-0	47
3			[(4-Hexylbenzene-1,3-diyl)bis(ox...	338	C18H34O2Si2	134273-01-1	47
4			9-Acridinol, trimethylsilyl ether	267	C16H17NOSi	1000463-65-5	46
5			Octopamine, 3TMS derivative	369	C17H35NO2Si3	068595-60-8	40



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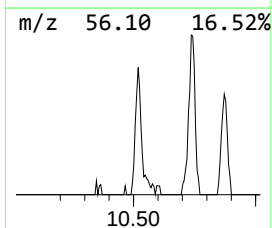
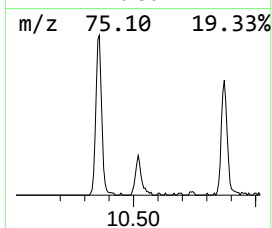
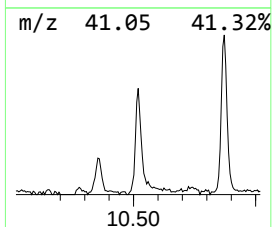
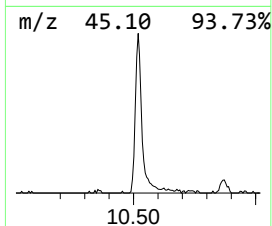
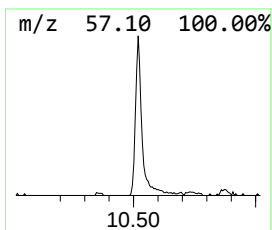
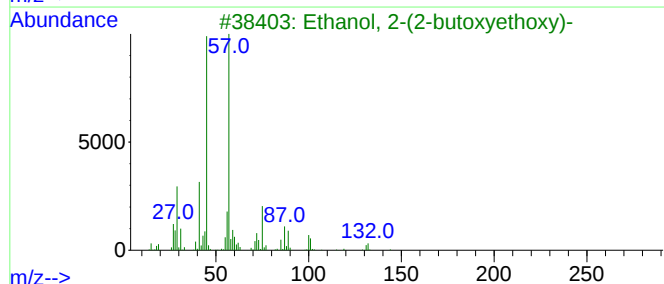
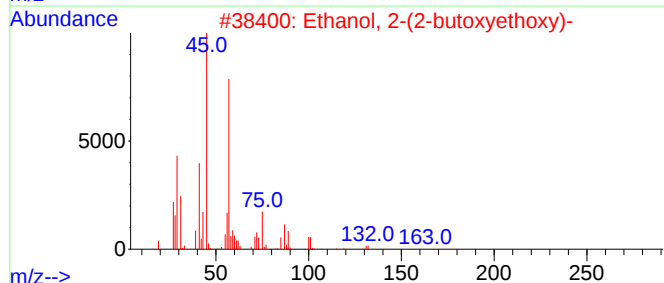
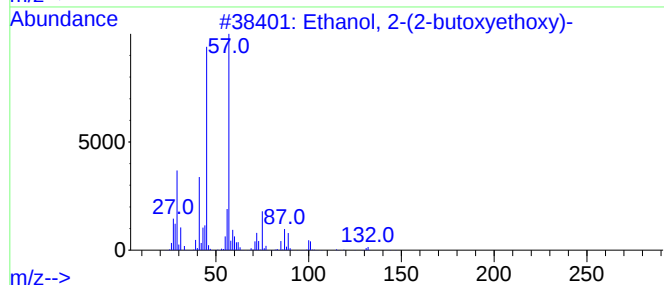
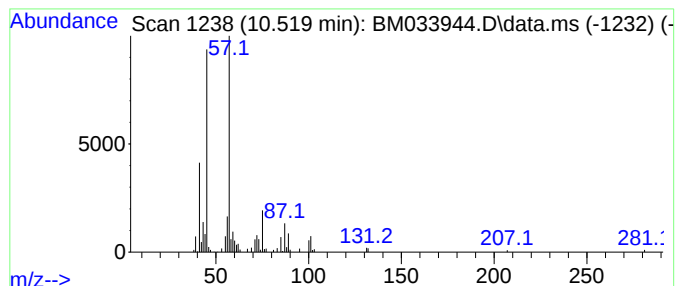
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.519	2.71 ng/ul	124650	Naphthalene-d8	10.742

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	86
3			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	83
4			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
5			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83



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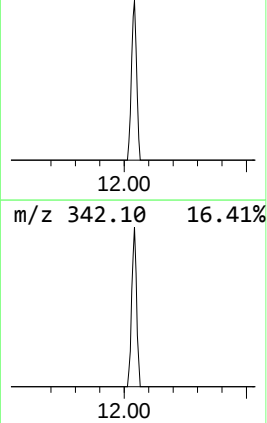
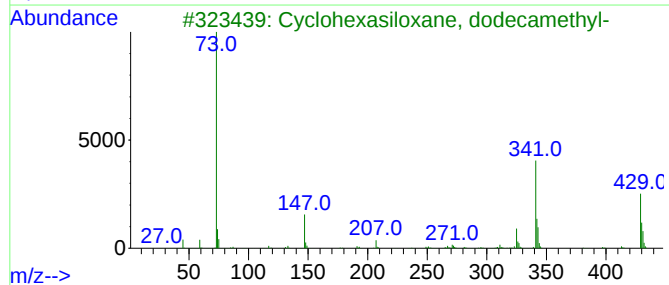
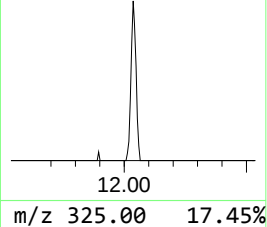
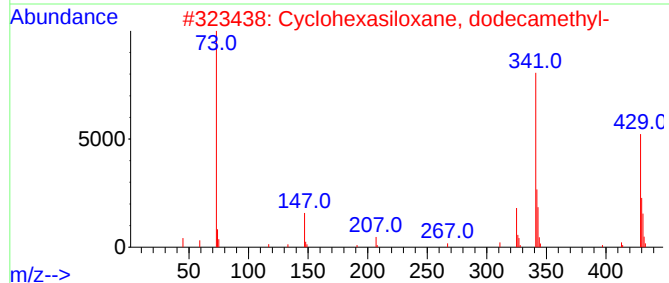
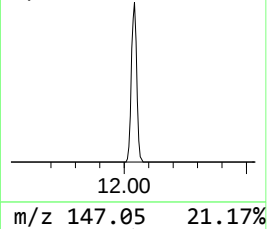
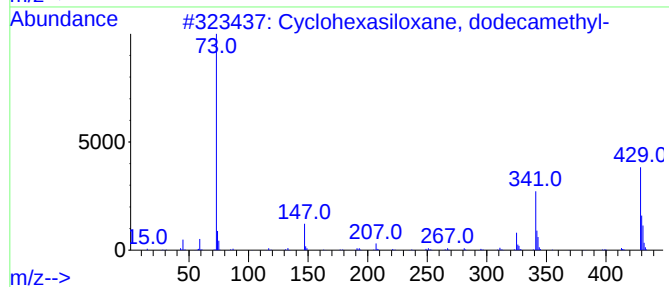
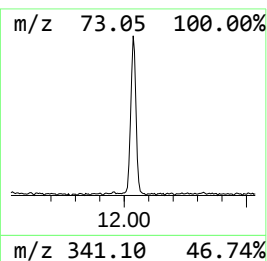
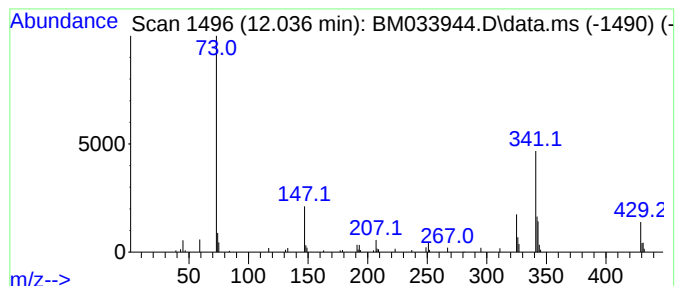
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Cyclohexasiloxane, dodecane... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.036	2.05 ng/ul	94162	Naphthalene-d8	10.742

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexasiloxane, dodecamethyl-	444	C12H36O6Si6	000540-97-6	91
2			Cyclohexasiloxane, dodecamethyl-	444	C12H36O6Si6	000540-97-6	91
3			Cyclohexasiloxane, dodecamethyl-	444	C12H36O6Si6	000540-97-6	83
4			4-Amino-1,2-benzenediol, O,O',N-...	341	C15H31N02Si3	1000493-87-4	50
5			4-Fluoro-N-(3-fluorophenyl)benze...	341	C15H17F2N02SSi	1000486-57-2	46



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ClientSampleId :
COHL9

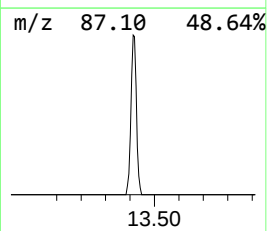
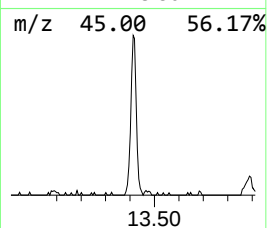
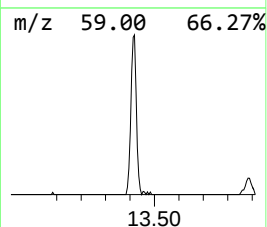
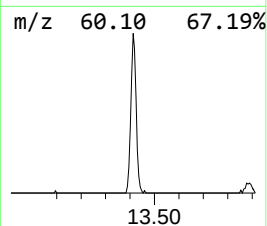
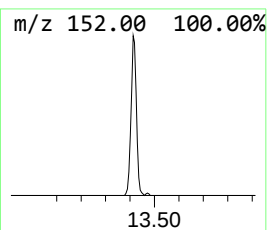
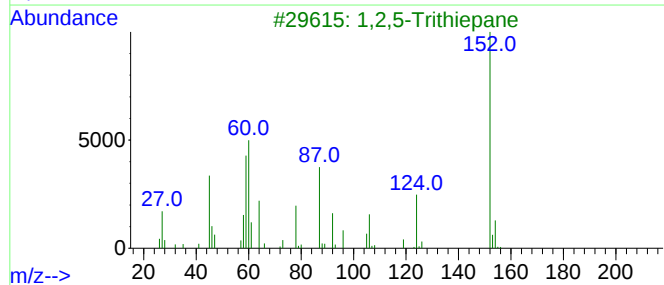
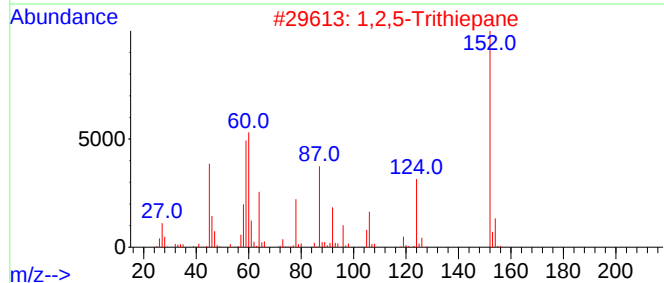
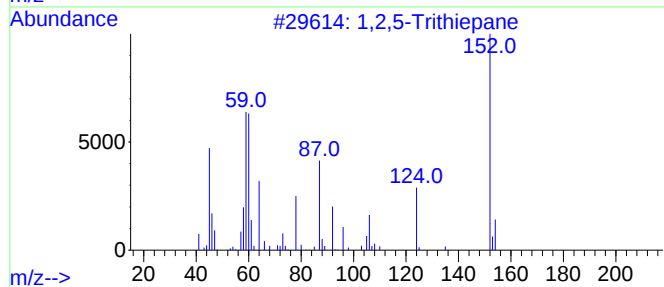
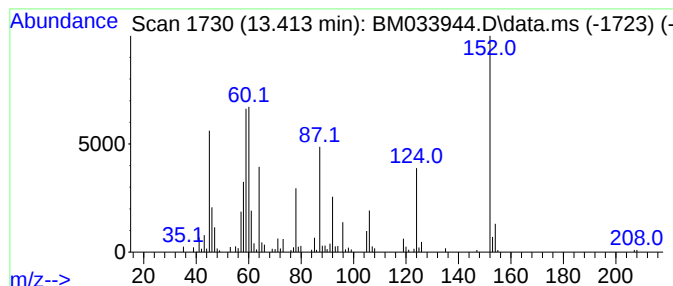
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,2,5-Trithiepane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.413	3.41 ng/ul	211192	Acenaphthene-d10	14.566

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,5-Trithiepane			152	C4H8S3	006576-93-8	99
2	1,2,5-Trithiepane			152	C4H8S3	006576-93-8	98
3	1,2,5-Trithiepane			152	C4H8S3	006576-93-8	97
4	1,2,4-Trithiolane, 3,5-dimethyl-			152	C4H8S3	023654-92-4	70
5	1,3,5-Trithiocycloheptane			152	C4H8S3	081451-19-6	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033944.D
Acq On : 01 Feb 2022 17:56
Operator : CG/JU
Sample : N1273-21
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
COHL9

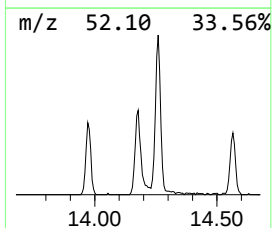
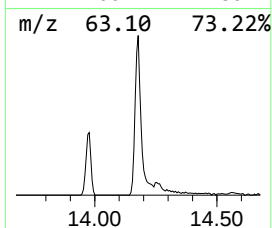
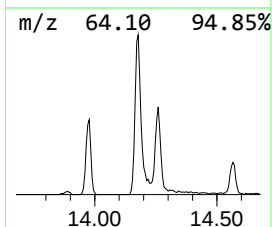
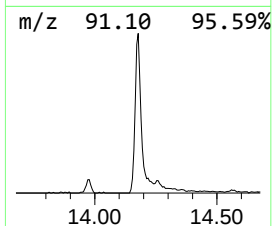
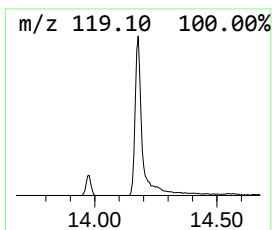
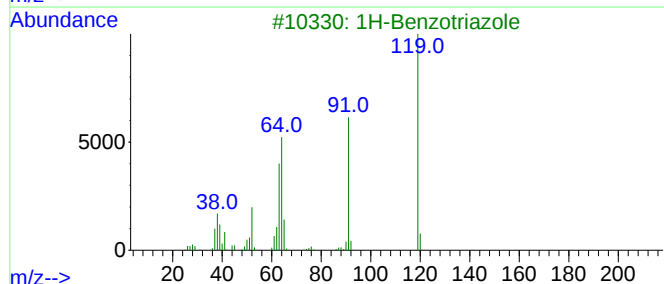
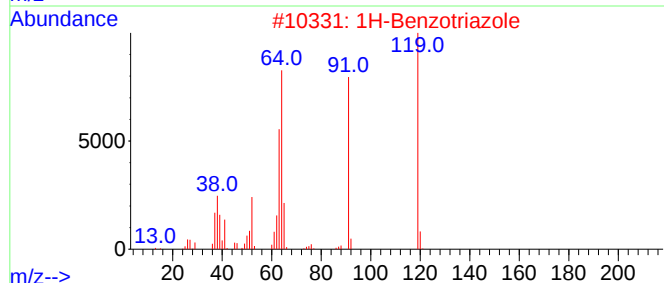
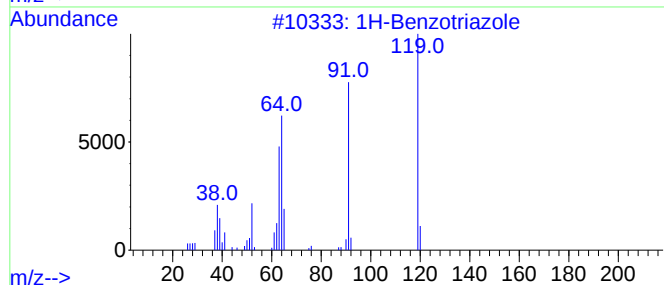
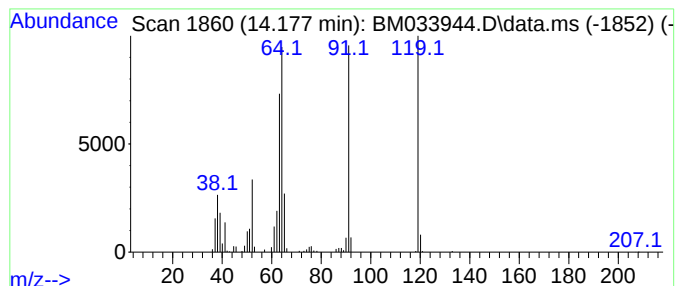
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 1H-Benzotriazole Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.177	6.71 ng/ul	415219	Acenaphthene-d10	14.566

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Benzotriazole			119	C6H5N3	000095-14-7	98
2	1H-Benzotriazole			119	C6H5N3	000095-14-7	98
3	1H-Benzotriazole			119	C6H5N3	000095-14-7	97
4	Ethane-1,2-diol, 1,2-bis(1-benzo...			296	C14H12N6O2	111507-88-1	91
5	1H-Benzotriazole			119	C6H5N3	000095-14-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
Data File : BM033944.D
Acq On : 01 Feb 2022 17:56
Operator : CG/JU
Sample : N1273-21
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0HL9

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.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
Cyclopentasilox...	9.542	3.0	ng/ul	136480	2	10.742	919380	20.0
Ethanol, 2-(2-b...	10.519	2.7	ng/ul	124650	2	10.742	919380	20.0
Cyclohexasiloxa...	12.036	2.0	ng/ul	94162	2	10.742	919380	20.0
1,2,5-Triethiepane	13.413	3.4	ng/ul	211192	3	14.566	1237680	20.0
1H-Benzotriazole	14.177	6.7	ng/ul	415219	3	14.566	1237680	20.0