

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
 Data File : BM036642.D
 Acq On : 22 Sep 2022 01:58
 Operator : CG/JU
 Sample : N4649-02
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EXZ18

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-BM091522.M
 Title : SVOA CALIBRATION

Signal : TIC: BM036642.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.122	3	6	28	rVB	6922816	9209301	78.56%	6.227%
2	3.393	47	52	60	rVB	84741	119728	1.02%	0.081%
3	3.822	121	125	138	rBV	438145	693967	5.92%	0.469%
4	4.063	161	166	171	rVB	46836	71036	0.61%	0.048%
5	4.157	177	182	186	rBV	29072	41631	0.36%	0.028%
6	4.704	270	275	279	rVV6	11150	19956	0.17%	0.013%
7	5.104	338	343	352	rVB	188370	272045	2.32%	0.184%
8	6.051	497	504	515	rVB	161529	259332	2.21%	0.175%
9	6.998	660	665	675	rBV	327463	621216	5.30%	0.420%
10	7.175	686	695	706	rVB	459313	804505	6.86%	0.544%
11	7.345	717	724	736	rVB	1819462	2827595	24.12%	1.912%
12	7.545	751	758	769	rBV	1628067	2599252	22.17%	1.758%
13	7.640	771	774	782	rVB6	10512	21657	0.18%	0.015%
14	7.757	789	794	799	rVB2	34257	54375	0.46%	0.037%
15	8.022	830	839	845	rBV	1500721	2398581	20.46%	1.622%
16	8.087	845	850	859	rVB	25131	46322	0.40%	0.031%
17	8.728	951	959	970	rBV	1344164	2153112	18.37%	1.456%
18	8.845	975	979	985	rVB	25875	42670	0.36%	0.029%
19	9.128	1021	1027	1030	rBV4	12120	22140	0.19%	0.015%
20	9.187	1030	1037	1044	rVB	2096641	3368668	28.74%	2.278%
21	9.910	1149	1160	1174	rBV	1596587	2662661	22.72%	1.801%
22	10.451	1244	1252	1264	rBV	2343731	3968740	33.86%	2.684%
23	10.616	1273	1280	1285	rVB5	11690	19650	0.17%	0.013%
24	10.833	1309	1317	1325	rBV	2215878	3710017	31.65%	2.509%
25	10.969	1332	1340	1358	rBV	2207889	3837032	32.73%	2.595%
26	11.122	1362	1366	1371	rVV2	11937	18245	0.16%	0.012%
27	11.186	1371	1377	1383	rVV	112794	195635	1.67%	0.132%
28	11.251	1383	1388	1392	rVV	25886	45157	0.39%	0.031%
29	11.310	1392	1398	1402	rVB3	12724	21219	0.18%	0.014%
30	11.369	1405	1408	1416	rVB8	13833	24285	0.21%	0.016%
31	11.486	1420	1428	1433	rBV	52919	95709	0.82%	0.065%
32	11.545	1433	1438	1447	rVB	131938	220668	1.88%	0.149%
33	12.033	1515	1521	1527	rBV	49007	78785	0.67%	0.053%
34	12.104	1527	1533	1536	rVV3	15490	33046	0.28%	0.022%
35	12.139	1536	1539	1548	rVB2	18564	32104	0.27%	0.022%
36	12.227	1548	1554	1560	rVB5	18964	30535	0.26%	0.021%

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
 Title : SVOA CALIBRATION

37	12.316	1564	1569	1574	rVB2	23613	34327	0.29%	0.023%
38	12.410	1579	1585	1588	rBV	51197	80329	0.69%	0.054%
39	12.451	1588	1592	1600	rVB	280981	420728	3.59%	0.285%
40	12.569	1605	1612	1617	rBV9	10528	24431	0.21%	0.017%

41	13.310	1733	1738	1746	rVB	51577	77442	0.66%	0.052%
42	13.551	1773	1779	1785	rBV3	22034	39190	0.33%	0.027%
43	13.957	1839	1848	1853	rBV5	18395	41816	0.36%	0.028%
44	14.057	1855	1865	1877	rBV	4766685	6475692	55.24%	4.379%
45	14.174	1882	1885	1892	rVB8	9998	21731	0.19%	0.015%

46	14.251	1892	1898	1900	rBV	714491	941912	8.04%	0.637%
47	14.280	1900	1903	1908	rVV	1357808	1805828	15.41%	1.221%
48	14.339	1908	1913	1926	rVB	5791350	8305609	70.86%	5.616%
49	14.469	1931	1935	1937	rVV5	11318	14690	0.13%	0.010%
50	14.498	1937	1940	1944	rVB	19904	26875	0.23%	0.018%

51	14.645	1958	1965	1973	rVB	3546886	5059220	43.16%	3.421%
52	14.751	1976	1983	1988	rVV2	50799	94463	0.81%	0.064%
53	14.833	1991	1997	2015	rVB	432900	700168	5.97%	0.473%
54	15.051	2030	2034	2039	rVB8	16884	27337	0.23%	0.018%
55	15.157	2046	2052	2062	rVB	114681	177026	1.51%	0.120%

56	15.327	2077	2081	2085	rVB4	14063	22451	0.19%	0.015%
57	15.439	2094	2100	2105	rBV7	15812	32927	0.28%	0.022%
58	15.633	2126	2133	2145	rBV	7324649	10058943	85.81%	6.802%
59	15.745	2145	2152	2163	rBV	1646713	2325084	19.84%	1.572%
60	15.868	2168	2173	2183	rVB	61325	112081	0.96%	0.076%

61	15.992	2189	2194	2200	rBV	74704	108616	0.93%	0.073%
62	16.110	2207	2214	2217	rVB	31099	42766	0.36%	0.029%
63	16.192	2222	2228	2234	rBV5	32596	55358	0.47%	0.037%
64	16.310	2243	2248	2251	rBV6	18538	31646	0.27%	0.021%
65	16.415	2262	2266	2269	rVB4	15340	20471	0.17%	0.014%

66	16.533	2279	2286	2302	rBV2	5203101	8120991	69.28%	5.492%
67	16.874	2340	2344	2347	rBV	70377	85691	0.73%	0.058%
68	16.915	2347	2351	2358	rVB	356128	438483	3.74%	0.297%
69	17.104	2379	2383	2386	rBV5	12664	20978	0.18%	0.014%
70	17.139	2386	2389	2392	rVV3	18285	25850	0.22%	0.017%

71	17.192	2396	2398	2403	rVB4	18274	21387	0.18%	0.014%
72	17.262	2406	2410	2415	rVB	58727	75999	0.65%	0.051%
73	17.380	2424	2430	2437	rVB	4337025	5772493	49.25%	3.903%
74	17.480	2440	2447	2457	rVV	7735295	10679070	91.10%	7.221%
75	17.945	2522	2526	2531	rBV	98535	125101	1.07%	0.085%

76	18.051	2541	2544	2545	rBV3	23025	22929	0.20%	0.016%
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77	18.204	2567	2570	2576	rVB	45915	61233	0.52%	0.041%
78	18.280	2578	2583	2591	rBV2	181191	395463	3.37%	0.267%
79	18.598	2634	2637	2648	rBV	190397	346099	2.95%	0.234%
80	18.886	2682	2686	2693	rVB	141391	183421	1.56%	0.124%
81	18.998	2699	2705	2718	rBV	7926520	9552362	81.49%	6.459%
82	19.180	2733	2736	2744	rVB2	117303	167619	1.43%	0.113%
83	19.327	2757	2761	2765	rBV2	93805	126359	1.08%	0.085%
84	19.398	2769	2773	2780	rVB	98585	145200	1.24%	0.098%
85	19.545	2794	2798	2804	rBV	269845	401773	3.43%	0.272%
86	19.680	2817	2821	2827	rBV	198621	284499	2.43%	0.192%
87	19.762	2829	2835	2845	rVV	9190431	11721951	100.00%	7.927%
88	19.956	2864	2868	2872	rBV	139461	180587	1.54%	0.122%
89	20.303	2925	2927	2932	rBV2	67900	77168	0.66%	0.052%
90	20.374	2936	2939	2945	rVB	275076	341630	2.91%	0.231%
91	20.639	2981	2984	2987	rBV	123846	140453	1.20%	0.095%
92	20.680	2988	2991	2997	rBV2	181958	216896	1.85%	0.147%
93	20.756	3001	3004	3007	rBV	177055	191467	1.63%	0.129%
94	21.003	3042	3046	3054	rBV	365513	474644	4.05%	0.321%
95	21.250	3085	3088	3093	rVB2	163692	283577	2.42%	0.192%
96	21.462	3121	3124	3132	rBV	234964	321979	2.75%	0.218%
97	21.545	3133	3138	3142	rBV	4208784	5392790	46.01%	3.647%
98	22.156	3239	3242	3250	rVB2	389447	698651	5.96%	0.472%
99	23.833	3520	3527	3534	rBV2	4393659	8378716	71.48%	5.666%
100	23.986	3547	3553	3563	rVB2	2270453	4589333	39.15%	3.103%

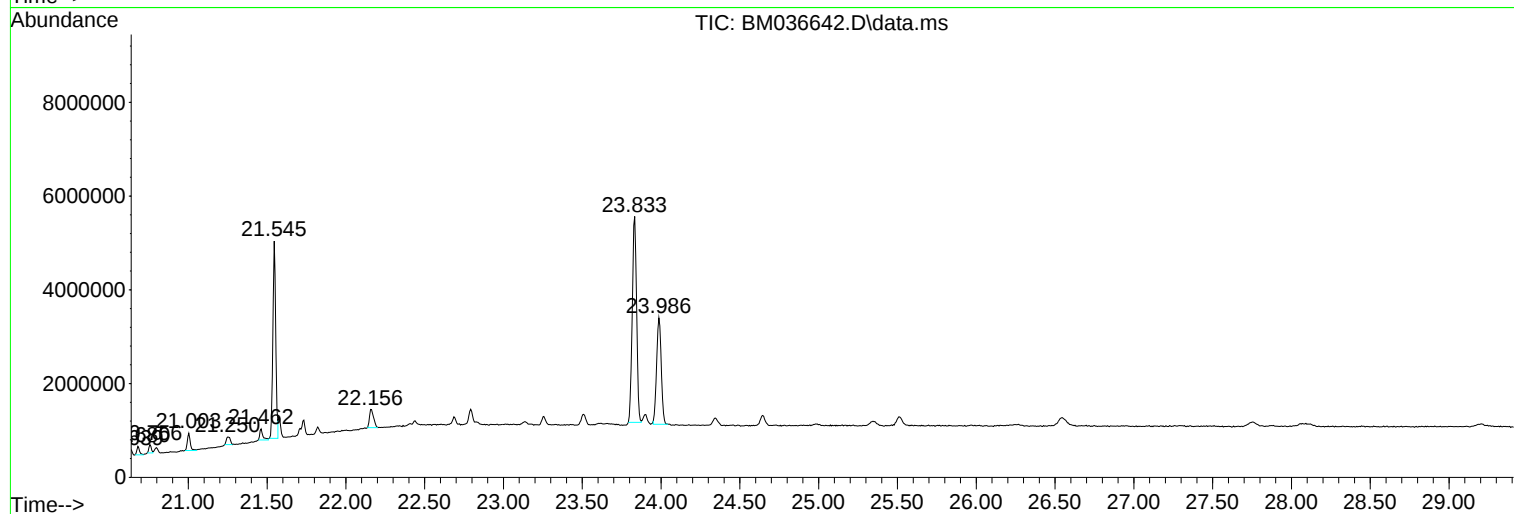
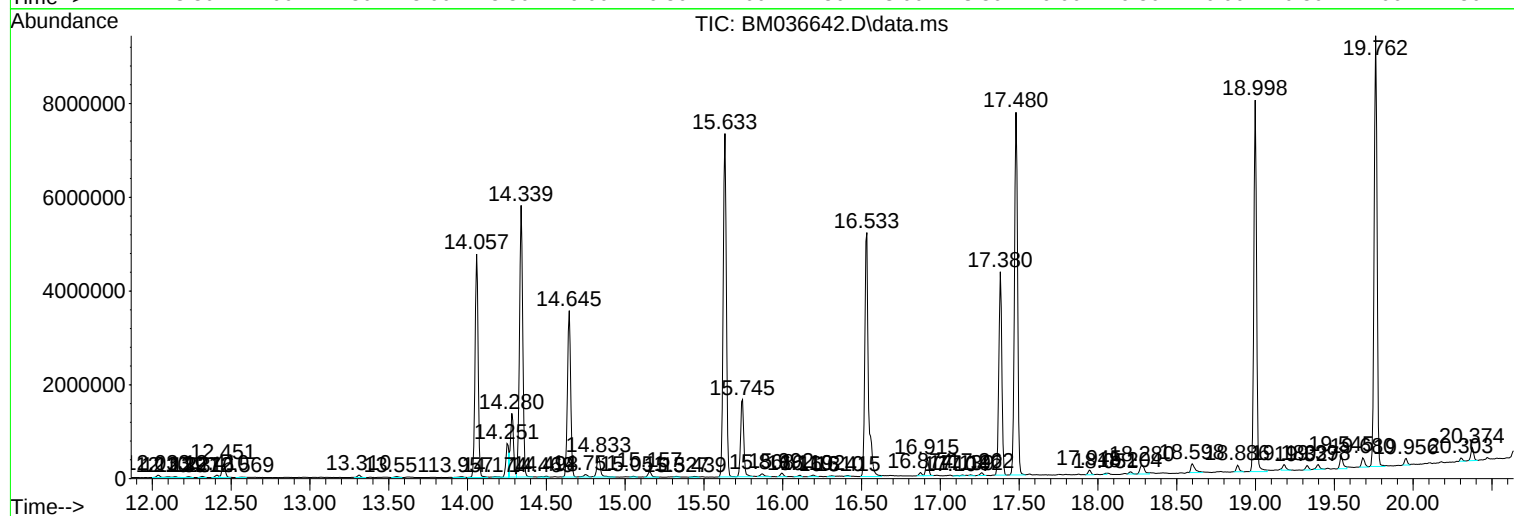
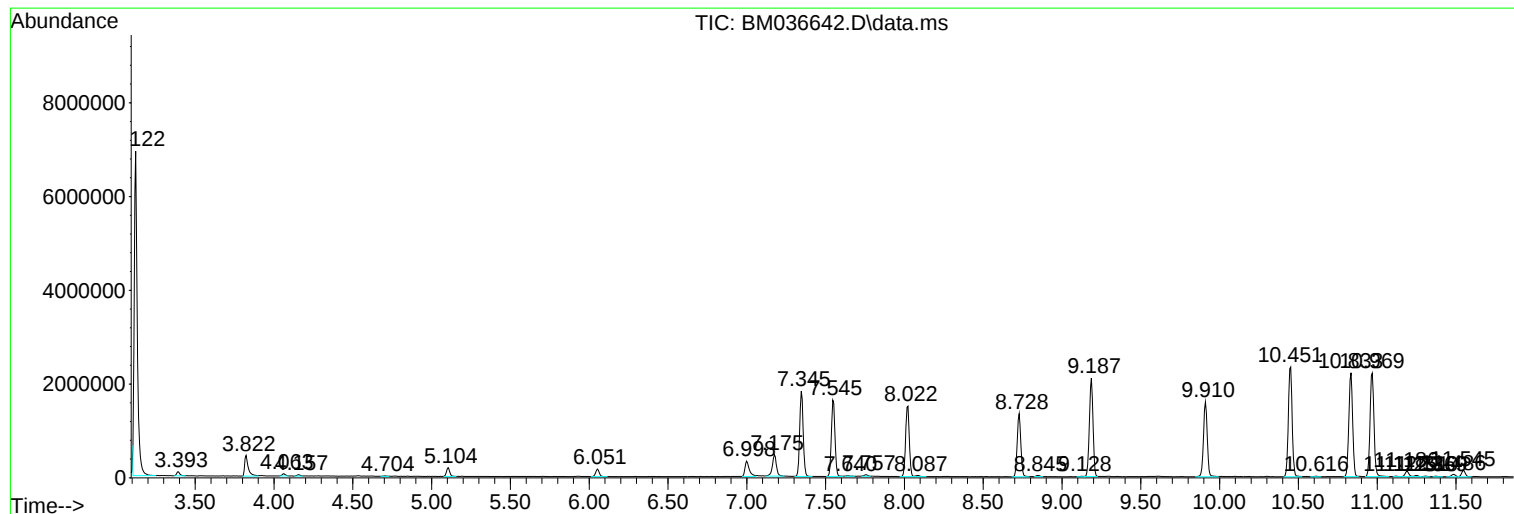
Sum of corrected areas: 147882576

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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



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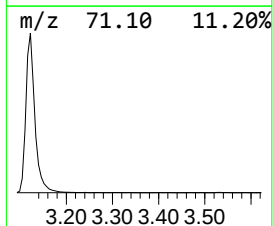
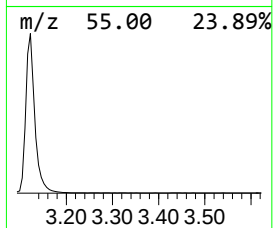
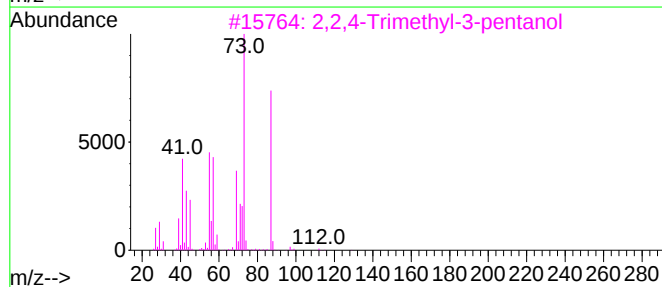
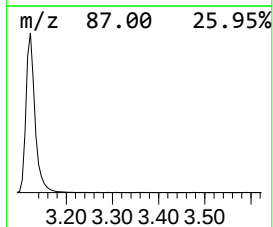
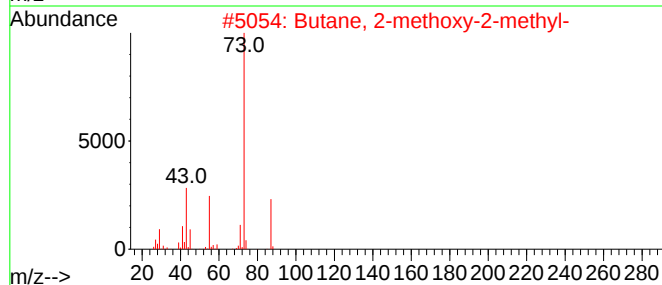
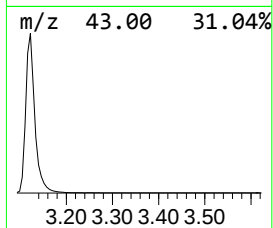
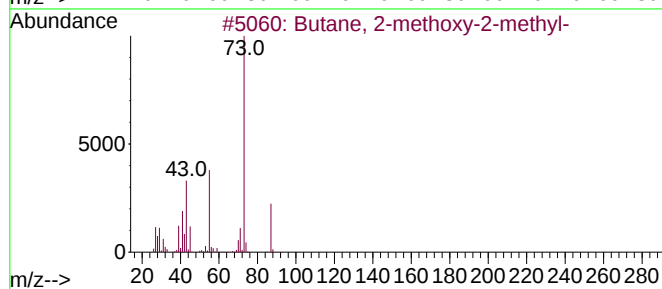
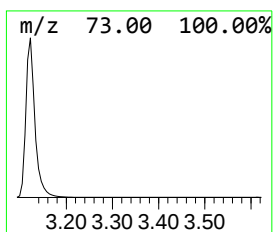
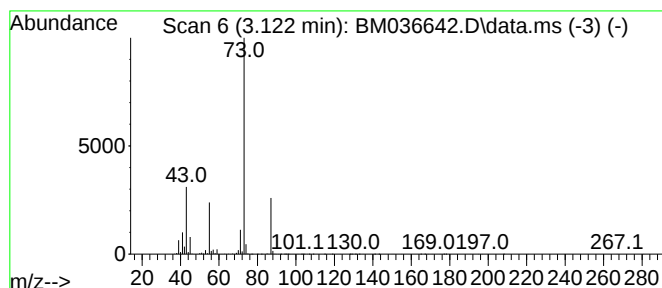
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.122	76.79 ng/ul	9209300	1,4-Dichlorobenzene-d4	8.022

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83		
2	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78		
3	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39		
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23		
5	Borane, dimethoxy-	74	C2H7BO2	004542-61-4	9		



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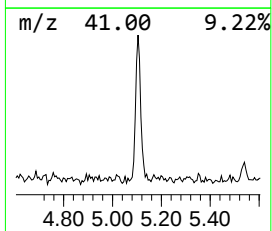
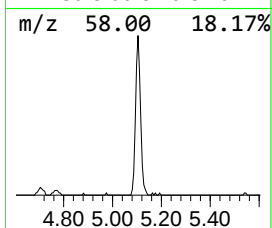
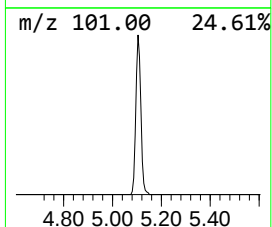
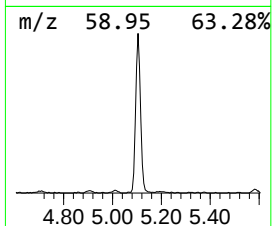
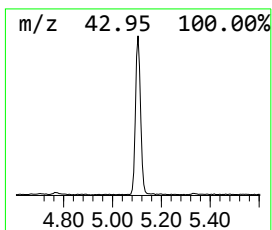
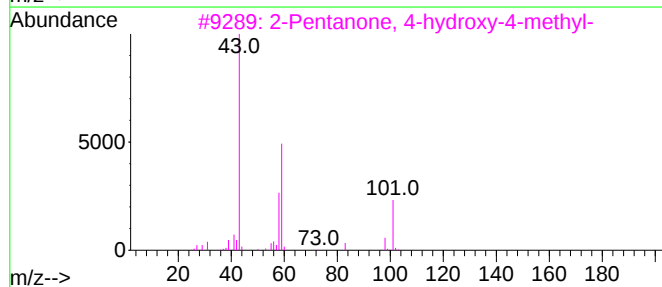
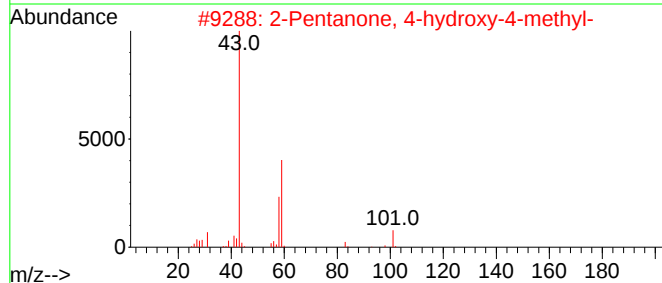
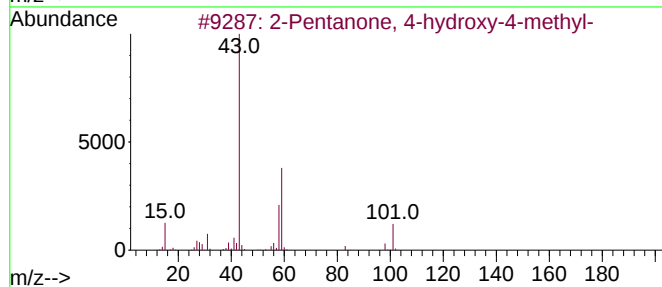
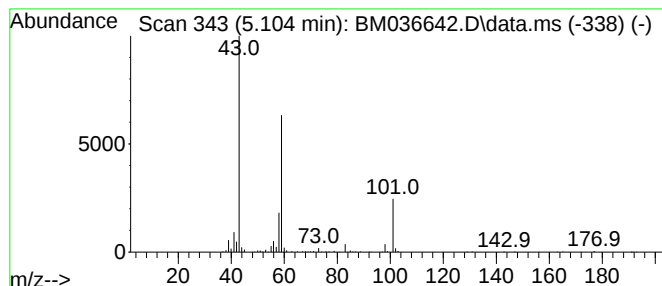
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.104	2.27 ng/ul	272045	1,4-Dichlorobenzene-d4	8.022

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
3			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	47
4			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
5			Tetraglyme	222	C10H22O5	000143-24-8	37



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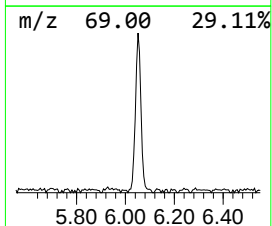
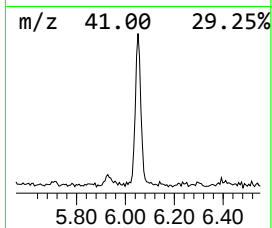
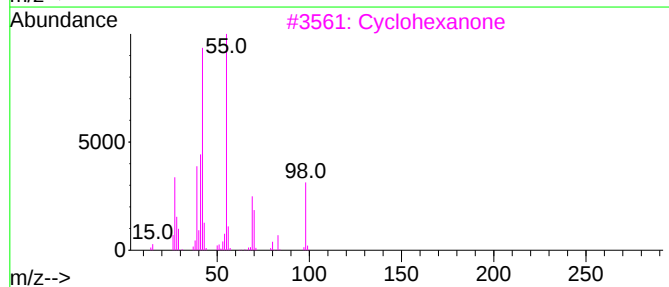
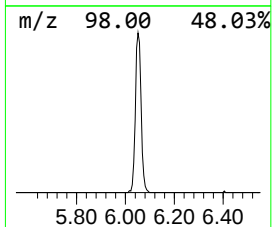
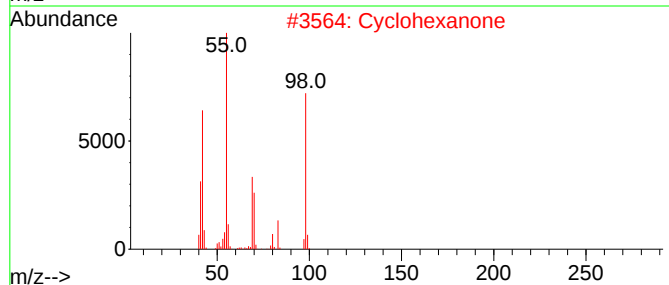
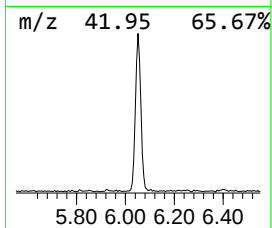
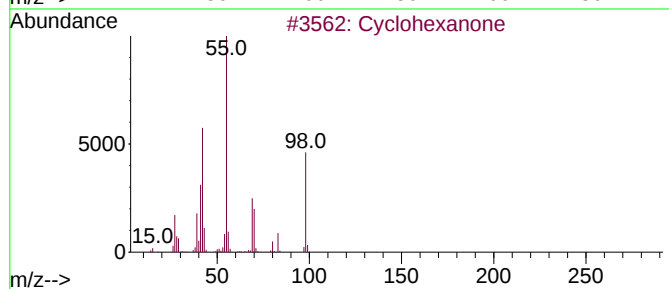
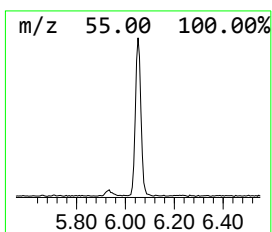
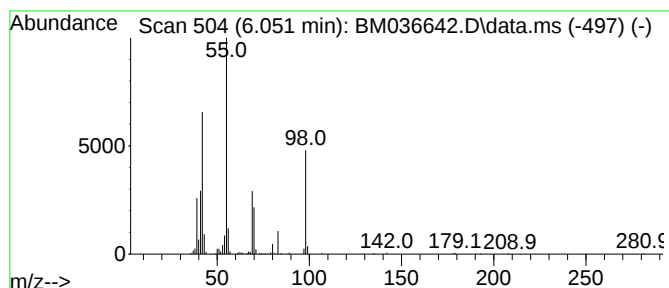
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Cyclohexanone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.051	2.16 ng/ul	259332	1,4-Dichlorobenzene-d4	8.022

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanone	98	C6H10O	000108-94-1	94
2			Cyclohexanone	98	C6H10O	000108-94-1	94
3			Cyclohexanone	98	C6H10O	000108-94-1	91
4			Cyclohexanone	98	C6H10O	000108-94-1	91
5			Cyclopentanone, 2-methyl-	98	C6H10O	001120-72-5	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
Data File : BM036642.D
Acq On : 22 Sep 2022 01:58
Operator : CG/JU
Sample : N4649-02
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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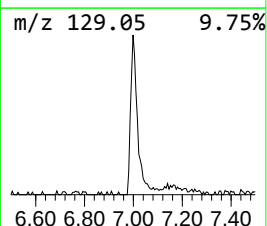
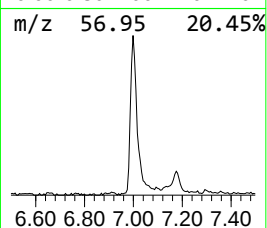
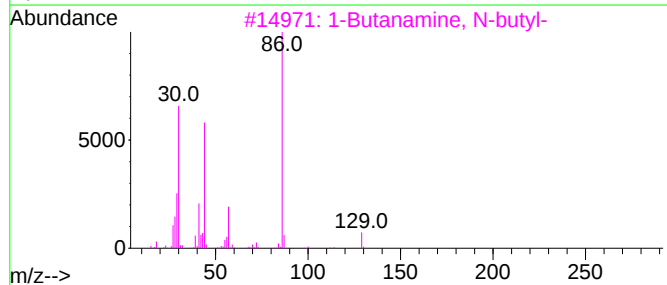
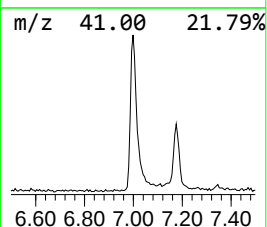
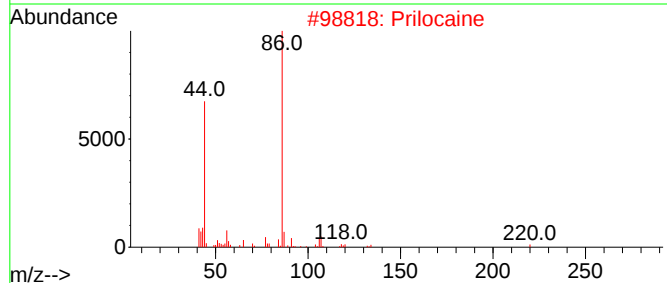
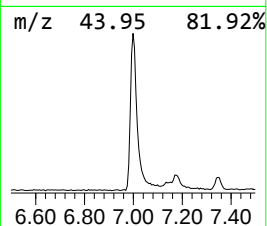
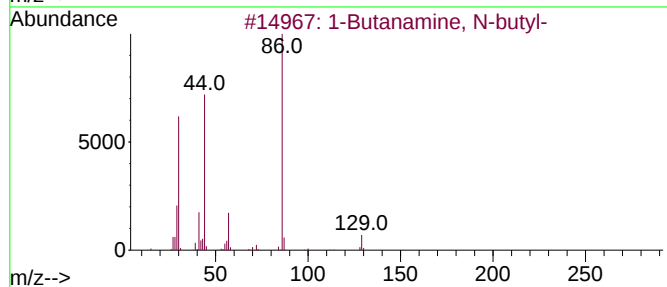
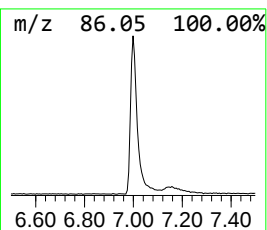
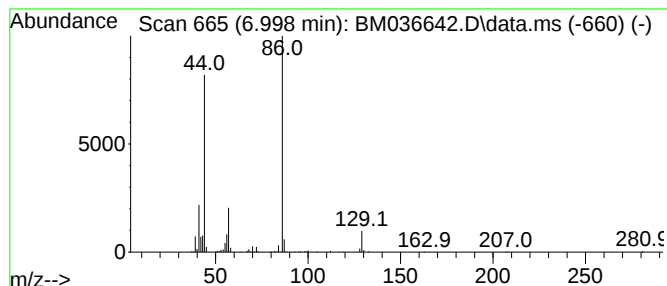
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Butanamine, N-butyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.998	5.18 ng/ul	621216	1,4-Dichlorobenzene-d4	8.022

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Butanamine, N-butyl-	129	C8H19N	000111-92-2	91
2			Prilocaine	220	C13H20N2O	000721-50-6	72
3			1-Butanamine, N-butyl-	129	C8H19N	000111-92-2	64
4			n-Butylethylenediamine	116	C6H16N2	019522-69-1	64
5			3',4'-Methylenedioxy-N-propylami...	235	C13H17NO3	1000457-81-1	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
Data File : BM036642.D
Acq On : 22 Sep 2022 01:58
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Sample : N4649-02
Misc :
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Instrument :
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ClientSampleId :
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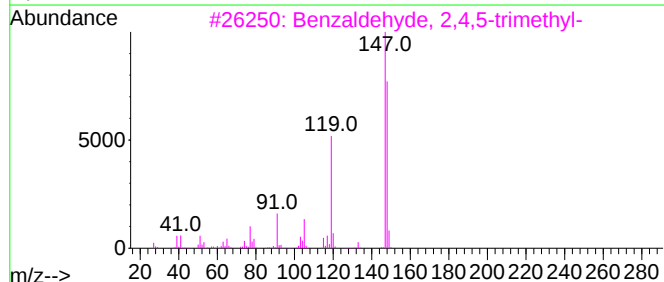
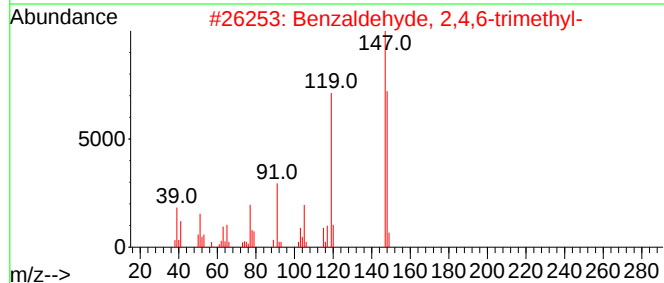
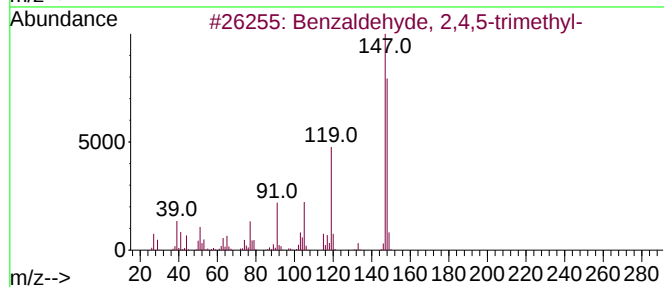
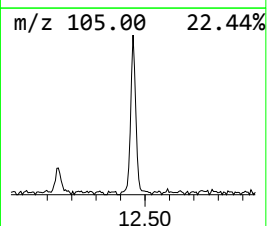
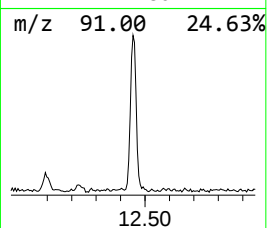
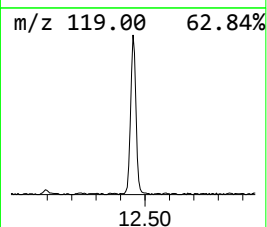
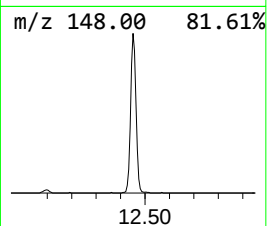
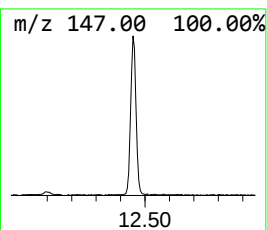
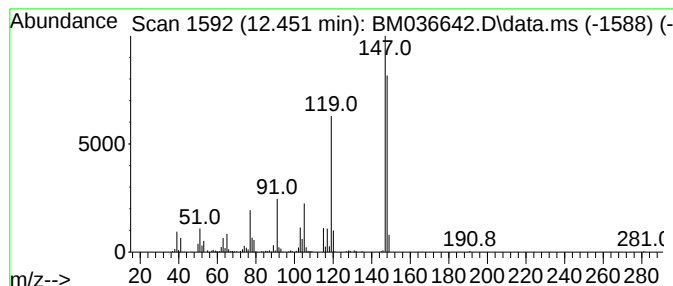
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzaldehyde, 2,4,5-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.451	2.27 ng/ul	420728	Naphthalene-d8	10.833

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzaldehyde, 2,4,5-trimethyl-	148	C10H12O	005779-72-6	95
2			Benzaldehyde, 2,4,6-trimethyl-	148	C10H12O	000487-68-3	95
3			Benzaldehyde, 2,4,5-trimethyl-	148	C10H12O	005779-72-6	95
4			Benzaldehyde, 2,4,6-trimethyl-	148	C10H12O	000487-68-3	94
5			Benzaldehyde, 2,4,6-trimethyl-	148	C10H12O	000487-68-3	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
Data File : BM036642.D
Acq On : 22 Sep 2022 01:58
Operator : CG/JU
Sample : N4649-02
Misc :
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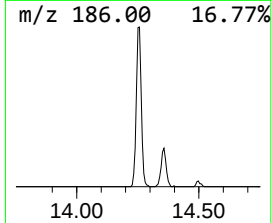
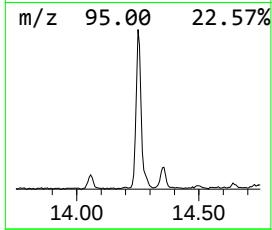
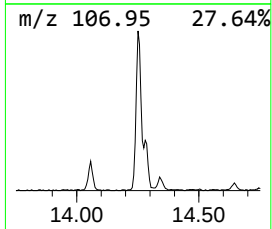
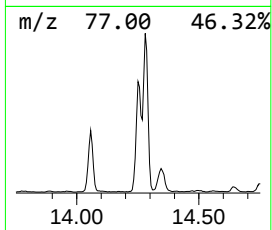
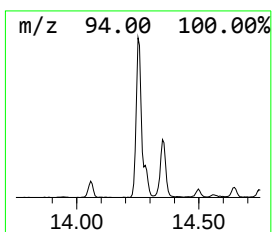
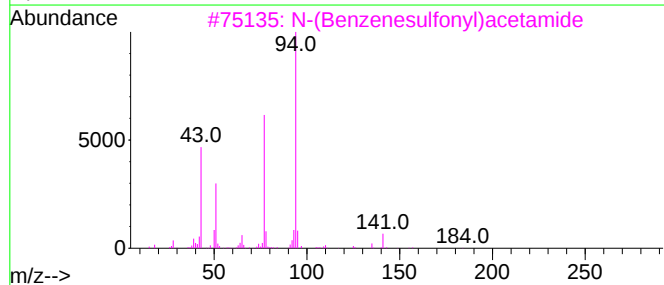
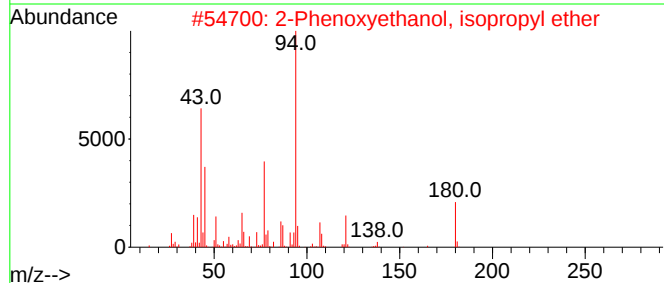
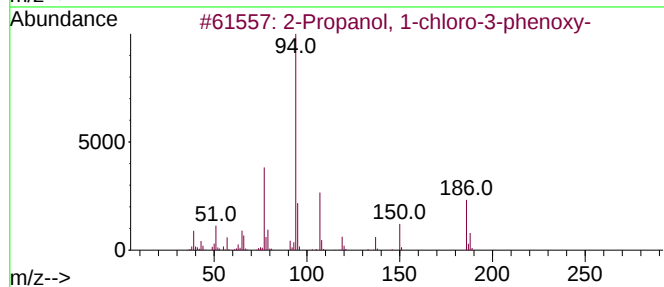
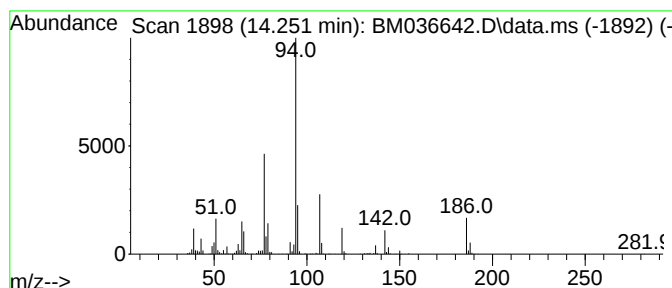
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 2-Propanol, 1-chloro-3-phen... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.251	3.72 ng/ul	941912	Acenaphthene-d10	14.645	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-chloro-3-phenoxy-	186	C9H11ClO2	004769-73-7	93
2	2-Phenoxyethanol, isopropyl ether	180	C11H16O2	1000394-79-3	53
3	N-(Benzenesulfonyl)acetamide	199	C8H9NO3S	005661-14-3	38
4	Oxirane, (phenoxymethyl)-	150	C9H10O2	000122-60-1	36
5	Carbonic acid, (3,4,4-trimethyl-...	252	C13H16O5	109123-69-5	36



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
Data File : BM036642.D
Acq On : 22 Sep 2022 01:58
Operator : CG/JU
Sample : N4649-02
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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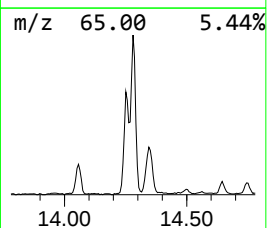
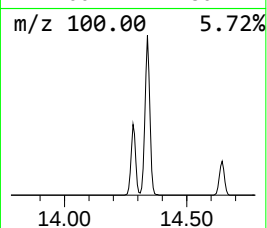
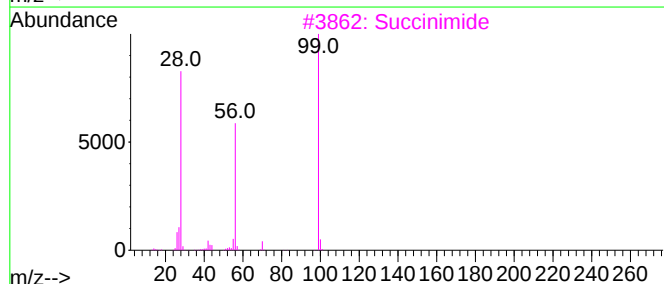
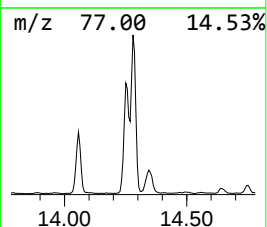
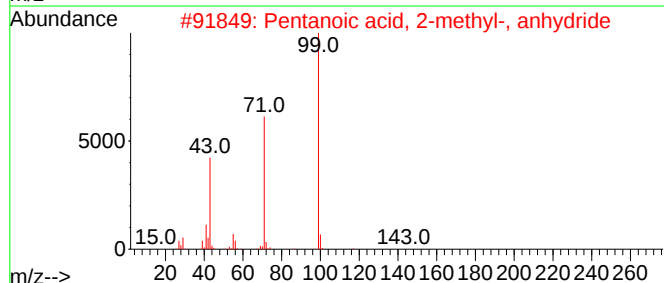
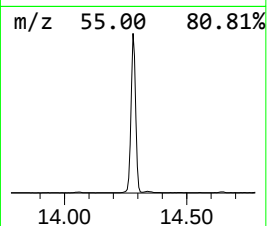
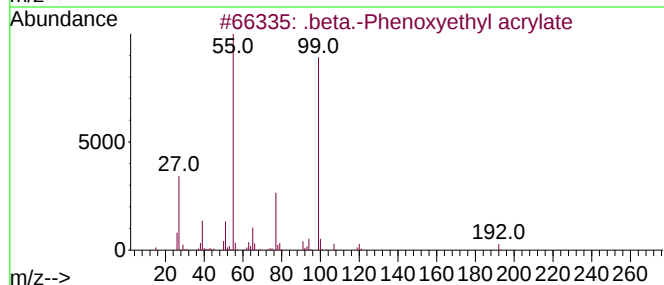
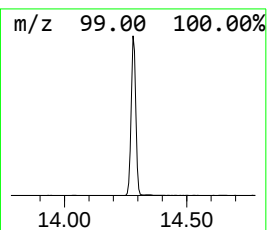
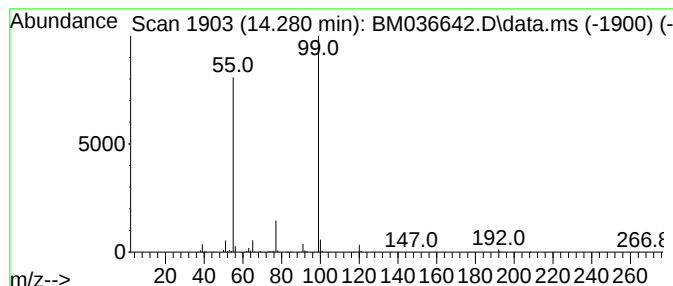
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.280	7.14 ng/ul	1805830	Acenaphthene-d10	14.645

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	.	beta.-Phenoxyethyl acrylate	192	C11H12O3	048145-04-6	9	
2	Pentanoic acid, 2-methyl-, anhyd...	214	C12H22O3	063169-61-9	9		
3	Succinimide	99	C4H5NO2	000123-56-8	5		
4	3,5-Diamino-1,2,4-triazole	99	C2H5N5	001455-77-2	5		
5	4-Methylthiazole	99	C4H5NS	000693-95-8	4		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
Data File : BM036642.D
Acq On : 22 Sep 2022 01:58
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Sample : N4649-02
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ALS Vial : 27 Sample Multiplier: 1

Instrument :
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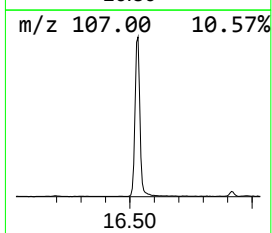
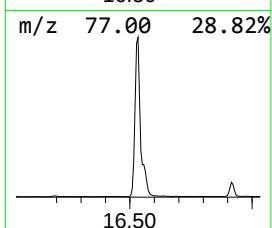
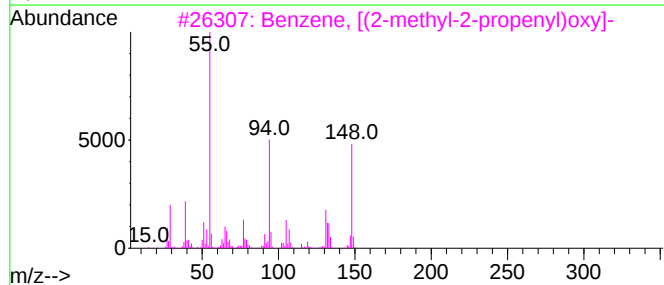
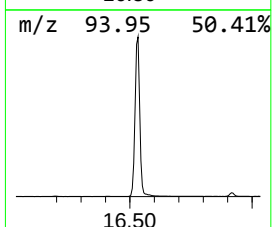
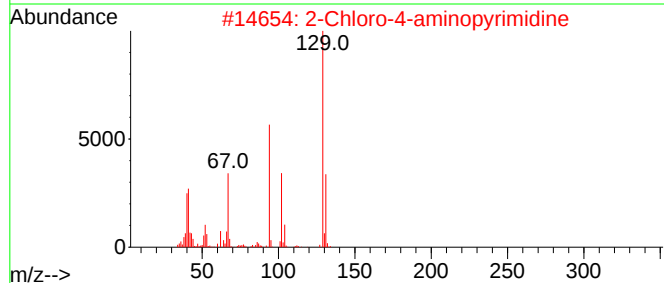
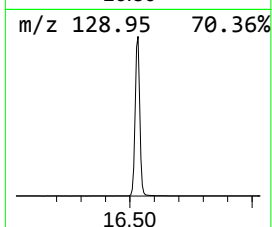
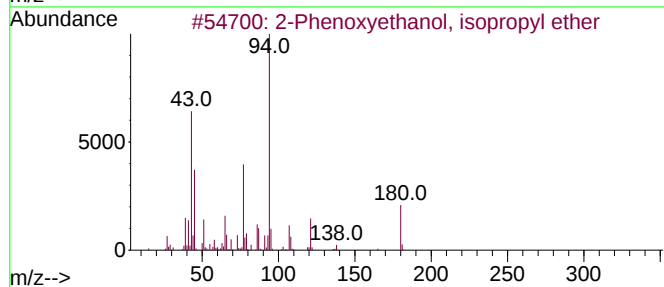
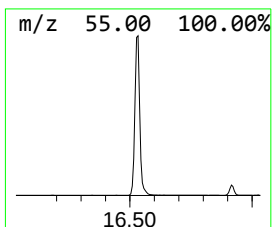
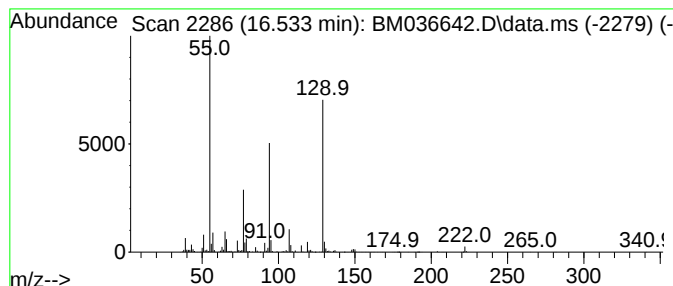
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-02 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.533	28.14 ng/ul	8120990	Phenanthrene-d10	17.380

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Phenoxyethanol, isopropyl ether	180	C11H16O2	1000394-79-3	38
2		2-Chloro-4-aminopyrimidine	129	C4H4ClN3	007461-50-9	37
3		Benzene, [(2-methyl-2-propenyl)oxy]-	148	C10H12O	005820-22-4	37
4		Oxirane, (phenoxymethyl)-	150	C9H10O2	000122-60-1	32
5		Methane, isopropoxyphenoxy-	166	C10H14O2	000828-12-6	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
Data File : BM036642.D
Acq On : 22 Sep 2022 01:58
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Sample : N4649-02
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ALS Vial : 27 Sample Multiplier: 1

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ClientSampleId :
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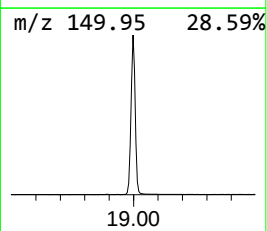
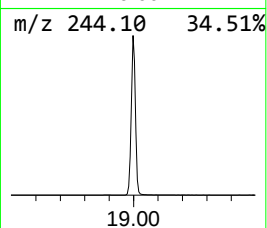
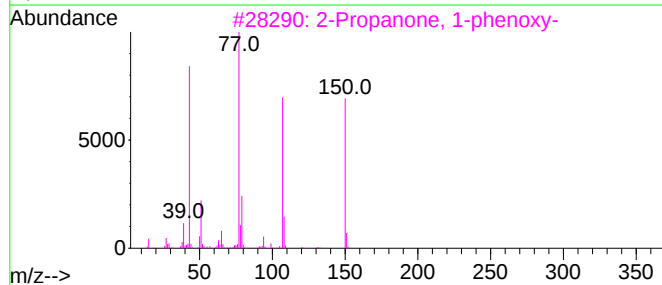
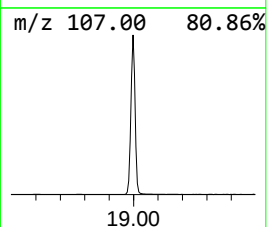
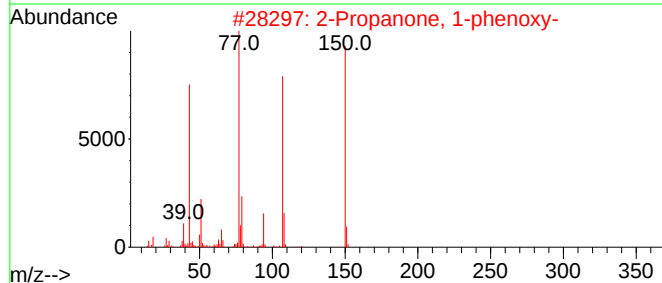
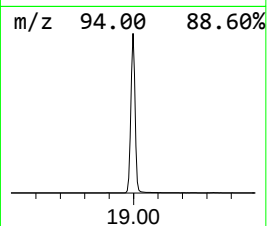
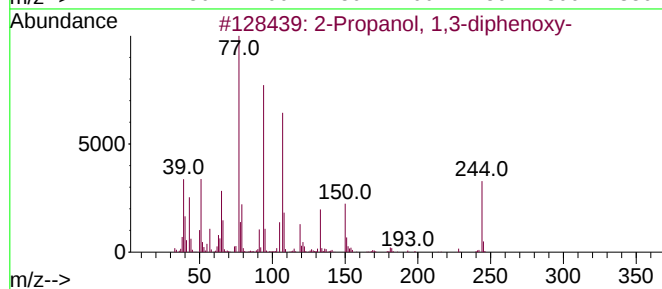
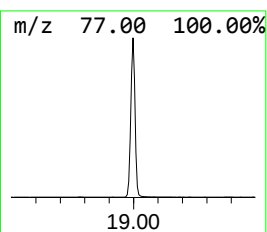
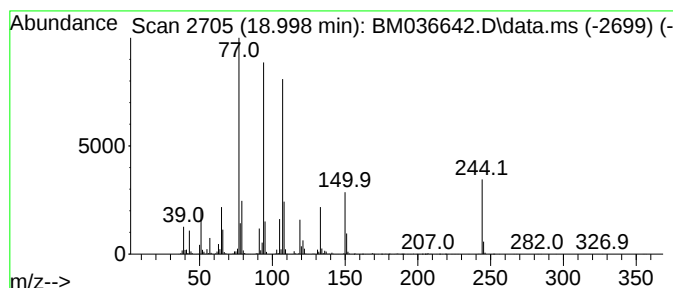
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 2-Propanol, 1,3-diphenoxy- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.998	33.10 ng/ul	9552360	Phenanthrene-d10	17.380

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1,3-diphenoxy-	244	C15H16O3		000622-04-8	98
2	2-Propanone, 1-phenoxy-	150	C9H10O2		000621-87-4	46
3	2-Propanone, 1-phenoxy-	150	C9H10O2		000621-87-4	43
4	Ethanol, 2-phenoxy-, benzoate	242	C15H14O3		004173-59-5	27
5	Phenol, 2-(2-methylpropyl)-	150	C10H14O		004167-75-3	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
Data File : BM036642.D
Acq On : 22 Sep 2022 01:58
Operator : CG/JU
Sample : N4649-02
Misc :
ALS Vial : 27 Sample Multiplier: 1

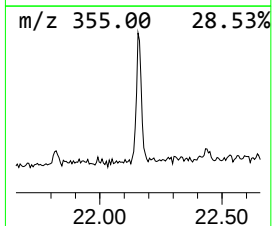
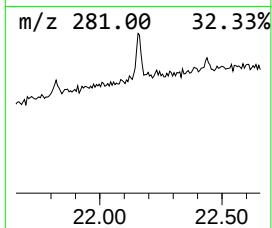
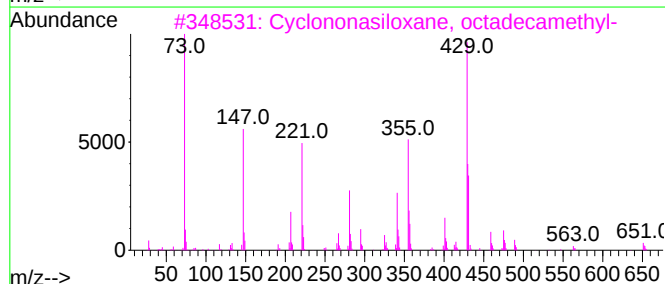
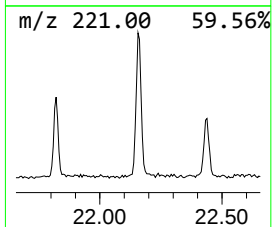
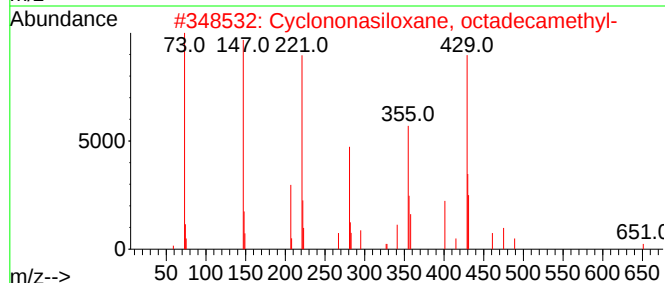
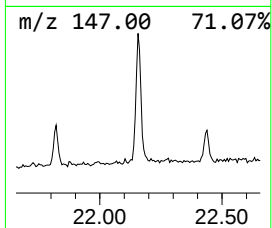
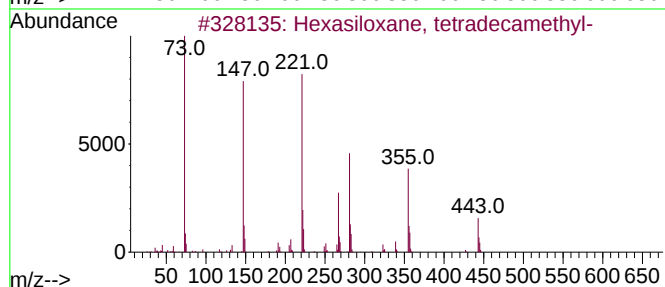
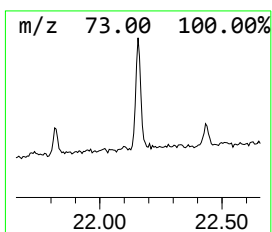
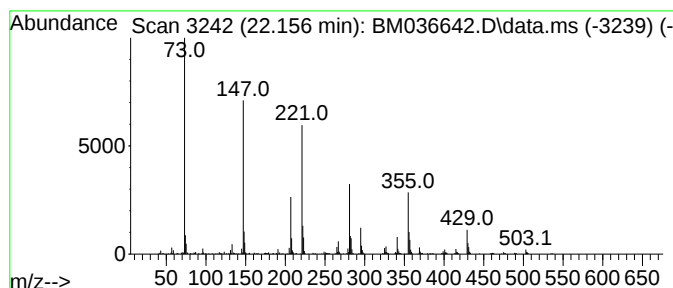
Instrument :
BNA_M
ClientSampleId :
EXZ18

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 Hexasiloxane, tetradecamethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD		R.T.	
22.156	2.59 ng/ul	698651	Chrysene-d12		21.544	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Hexasiloxane, tetradecamethyl-		458	C14H42O5Si6	000107-52-8	74
2	Cyclononasiloxane, octadecamethyl-		666	C18H54O9Si9	000556-71-8	64
3	Cyclononasiloxane, octadecamethyl-		666	C18H54O9Si9	000556-71-8	50
4	3,6-Dioxo-2,4,5,7-tetrasilaoctan...		294	C10H30O2Si4	004342-25-0	38
5	1,1,1,5,7,7,7-Heptamethyl-3,3-bi...		444	C13H40O5Si6	038147-00-1	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092122\
Data File : BM036642.D
Acq On : 22 Sep 2022 01:58
Operator : CG/JU
Sample : N4649-02
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EXZ18

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.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
Butane, 2-metho...	3.122	76.8	ng/ul	9209300	1	8.022	2398580	20.0
2-Pentanone, 4-...	5.104	2.3	ng/ul	272045	1	8.022	2398580	20.0
Cyclohexanone	6.051	2.2	ng/ul	259332	1	8.022	2398580	20.0
1-Butanamine, N...	6.998	5.2	ng/ul	621216	1	8.022	2398580	20.0
Benzaldehyde, 2...	12.451	2.3	ng/ul	420728	2	10.833	3710020	20.0
2-Propanol, 1-C...	14.251	3.7	ng/ul	941912	3	14.645	5059220	20.0
unknown-01	14.280	7.1	ng/ul	1805830	3	14.645	5059220	20.0
unknown-02	16.533	28.1	ng/ul	8120990	4	17.380	5772490	20.0
2-Propanol, 1,3...	18.998	33.1	ng/ul	9552360	4	17.380	5772490	20.0
Hexasiloxane, t...	22.156	2.6	ng/ul	698651	5	21.544	5392790	20.0