

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112018\
 Data File : BM017703.D
 Acq On : 20 Nov 2018 14:11
 Operator : MJ/SJ
 Sample : J6001-01
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A40T1

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111918MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.187	31	34	39	rBV	58769	77873	0.61%	0.034%
2	3.228	39	41	49	rVB5	14811	23237	0.18%	0.010%
3	3.505	82	88	92	rBV3	18281	34588	0.27%	0.015%
4	3.887	149	153	155	rBV3	18276	24191	0.19%	0.010%
5	4.152	192	198	202	rBV5	22671	42008	0.33%	0.018%
6	4.199	202	206	214	rVB	46863	77018	0.61%	0.033%
7	4.781	299	305	315	rBV	207450	318940	2.51%	0.137%
8	4.969	330	337	346	rBV	99975	153081	1.20%	0.066%
9	5.828	478	483	493	rVB	91857	147898	1.16%	0.064%
10	6.022	512	516	521	rVB3	10592	15455	0.12%	0.007%
11	6.287	559	561	570	rVB3	11850	22053	0.17%	0.009%
12	6.498	592	597	601	rBV3	12981	18460	0.15%	0.008%
13	6.781	636	645	660	rBV	1178289	2058163	16.20%	0.887%
14	6.946	666	673	685	rBV	927459	1441060	11.34%	0.621%
15	7.134	698	705	708	rBV	1249872	2108745	16.59%	0.908%
16	7.163	708	710	722	rVV	1275826	1849270	14.55%	0.797%
17	7.251	722	725	733	rVB8	8526	17023	0.13%	0.007%
18	7.598	773	784	796	rBV	750611	1188370	9.35%	0.512%
19	8.128	864	874	880	rBV2	1080677	2696326	21.22%	1.161%
20	8.175	880	882	890	rVB	117555	173025	1.36%	0.075%
21	8.310	898	905	916	rBV	1476650	2473220	19.46%	1.065%
22	8.422	916	924	936	rVB	1604363	2642320	20.79%	1.138%
23	8.657	957	964	970	rVB	586246	934555	7.35%	0.403%
24	8.751	973	980	983	rBV	1205377	1917362	15.09%	0.826%
25	8.792	983	987	996	rVB	1812028	2970930	23.38%	1.280%
26	9.051	1024	1031	1042	rBV	2389449	3640567	28.65%	1.568%
27	9.463	1095	1101	1104	rBV	1031056	1842530	14.50%	0.794%
28	9.498	1104	1107	1113	rVB	1260223	1865829	14.68%	0.804%
29	9.998	1182	1192	1215	rBV2	1459357	4292466	33.78%	1.849%
30	10.304	1238	1244	1248	rBV2	56875	109094	0.86%	0.047%
31	10.369	1248	1255	1259	rVV	1078400	1870812	14.72%	0.806%
32	10.416	1259	1263	1274	rVB	917821	1475191	11.61%	0.635%
33	10.528	1274	1282	1299	rBV	432343	985710	7.76%	0.425%
34	11.081	1369	1376	1382	rBV	35841	55381	0.44%	0.024%

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35	11.669	1468	1476	1495	rBV	2072460	3685225	29.00%	1.587%
36	11.910	1510	1517	1533	rBV	1879124	3284256	25.84%	1.415%
37	12.263	1565	1577	1588	rBV	3233155	5245855	41.28%	2.260%
38	12.410	1596	1602	1615	rVB	1145054	1785941	14.05%	0.769%
39	12.663	1639	1645	1653	rBV	1624083	2516334	19.80%	1.084%
40	12.733	1653	1657	1666	rVB	59673	106439	0.84%	0.046%
41	13.069	1707	1714	1722	rBV	3567192	5105757	40.18%	2.199%
42	13.333	1751	1759	1772	rBV	2152453	3370785	26.52%	1.452%
43	13.433	1772	1776	1790	rVB2	54714	89210	0.70%	0.038%
44	13.663	1809	1815	1823	rBV	2902160	4149108	32.65%	1.787%
45	13.739	1823	1828	1838	rVV	720742	1064177	8.37%	0.458%
46	13.833	1838	1844	1854	rVV	1865215	2830933	22.28%	1.219%
47	13.933	1854	1861	1863	rVV	3339506	4815959	37.90%	2.074%
48	13.963	1863	1866	1880	rVB	4481529	6260987	49.27%	2.697%
49	14.245	1906	1914	1920	rBV2	1655878	2550880	20.07%	1.099%
50	14.310	1920	1925	1933	rVB	418360	609696	4.80%	0.263%
51	14.486	1945	1955	1973	rBV3	2688248	5764932	45.36%	2.483%
52	14.639	1973	1981	1991	rVV2	3124350	6049470	47.60%	2.606%
53	14.710	1991	1993	2002	rVB7	20978	30846	0.24%	0.013%
54	14.904	2019	2026	2033	rBV	560221	777704	6.12%	0.335%
55	15.092	2051	2058	2063	rBV	4067164	5541701	43.61%	2.387%
56	15.245	2077	2084	2089	rBV	4376017	6084406	47.88%	2.621%
57	15.298	2089	2093	2101	rVB	882190	1214195	9.55%	0.523%
58	15.380	2101	2107	2119	rBV	1589061	2147403	16.90%	0.925%
59	15.521	2120	2131	2138	rBV	5334671	7640926	60.13%	3.291%
60	15.763	2165	2172	2177	rBV	179051	309192	2.43%	0.133%
61	15.886	2189	2193	2201	rVB	124938	174769	1.38%	0.075%
62	16.121	2229	2233	2237	rBV	23114	28216	0.22%	0.012%
63	16.192	2237	2245	2252	rVV3	75903	167788	1.32%	0.072%
64	16.298	2257	2263	2275	rVB2	3755423	5856627	46.09%	2.523%
65	16.651	2317	2323	2336	rBV	2579074	3701394	29.13%	1.594%
66	16.786	2341	2346	2349	rVV4	19120	32881	0.26%	0.014%
67	16.821	2349	2352	2357	rVB4	11811	17239	0.14%	0.007%
68	16.998	2376	2382	2386	rBV	2257405	3297464	25.95%	1.420%
69	17.039	2386	2389	2394	rVV	3942692	5435976	42.78%	2.341%
70	17.098	2394	2399	2408	rVV	4978002	6719358	52.87%	2.894%
71	17.168	2408	2411	2420	rVB8	20746	46960	0.37%	0.020%

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72	17.345	2437	2441	2444	rBV5	11352	15155	0.12%	0.007%
73	17.621	2484	2488	2491	rBV3	15198	20113	0.16%	0.009%
74	17.910	2534	2537	2543	rBV6	30077	49510	0.39%	0.021%
75	17.980	2543	2549	2559	rVV	4875985	5957121	46.88%	2.566%
76	18.257	2592	2596	2601	rVB4	21834	28831	0.23%	0.012%
77	19.039	2724	2729	2733	rBV	52343	70962	0.56%	0.031%
78	19.198	2752	2756	2762	rBV6	28148	44516	0.35%	0.019%
79	19.286	2766	2771	2779	rBV	718746	1043860	8.21%	0.450%
80	19.404	2784	2791	2794	rVV2	5405301	8489759	66.81%	3.657%
81	19.427	2794	2795	2811	rVB	2469761	2690830	21.17%	1.159%
82	19.651	2828	2833	2841	rBV	289928	392574	3.09%	0.169%
83	19.809	2856	2860	2865	rVB3	38542	54279	0.43%	0.023%
84	20.274	2936	2939	2942	rBV4	16918	22263	0.18%	0.010%
85	20.451	2964	2969	2977	rBV	580402	742571	5.84%	0.320%
86	20.556	2984	2987	2992	rBV5	29892	50052	0.39%	0.022%
87	21.133	3081	3085	3089	rBV	199279	289287	2.28%	0.125%
88	21.186	3089	3094	3105	rVB2	7126583	12708086	100.00%	5.474%
89	21.474	3139	3143	3151	rBV2	464876	626214	4.93%	0.270%
90	21.633	3164	3170	3175	rBV	8563443	9633743	75.81%	4.150%
91	21.750	3186	3190	3194	rBV2	270629	398875	3.14%	0.172%
92	21.997	3227	3232	3237	rBV	4177968	5201618	40.93%	2.240%
93	22.239	3271	3273	3279	rVB	134276	156062	1.23%	0.067%
94	22.739	3351	3358	3368	rVB	4362547	7038333	55.38%	3.032%
95	23.256	3438	3446	3450	rBV	4053082	8373430	65.89%	3.607%
96	23.297	3450	3453	3462	rVV	2350077	3649064	28.71%	1.572%
97	23.392	3462	3469	3478	rVB	1830741	3545184	27.90%	1.527%
98	23.903	3552	3556	3564	rVB2	232466	426316	3.35%	0.184%
99	25.580	3831	3841	3855	rVB	3119901	8363339	65.81%	3.602%
100	26.227	3942	3951	3967	rVB	1374829	4008976	31.55%	1.727%

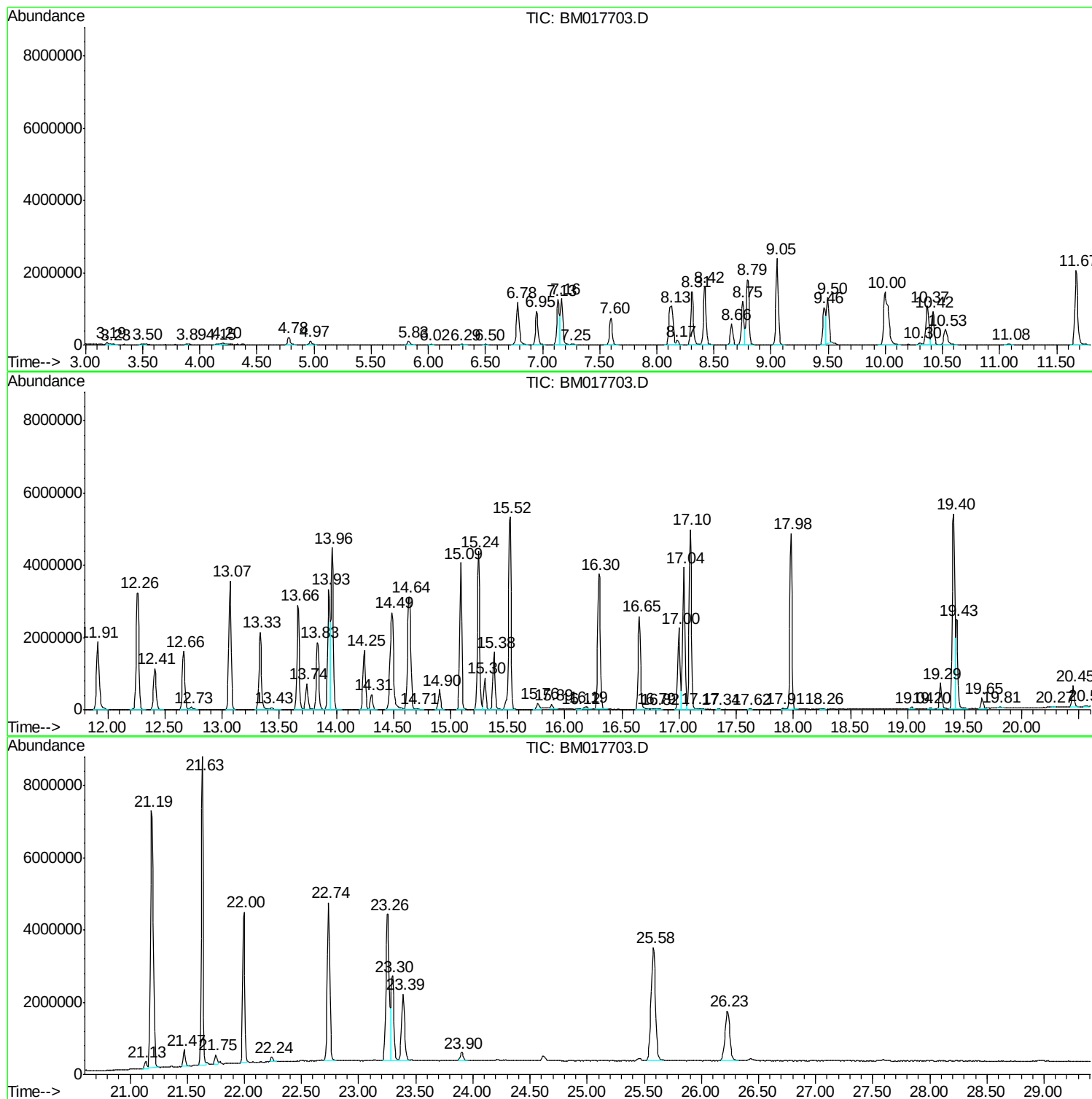
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111918MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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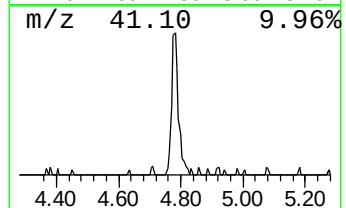
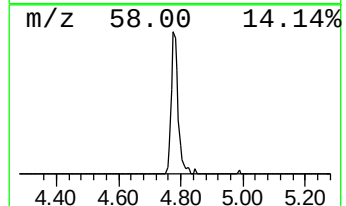
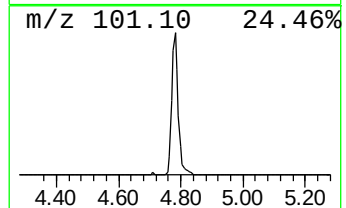
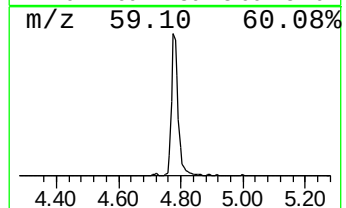
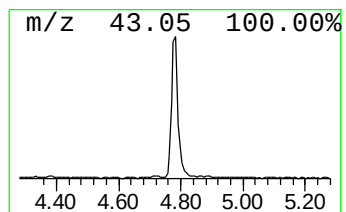
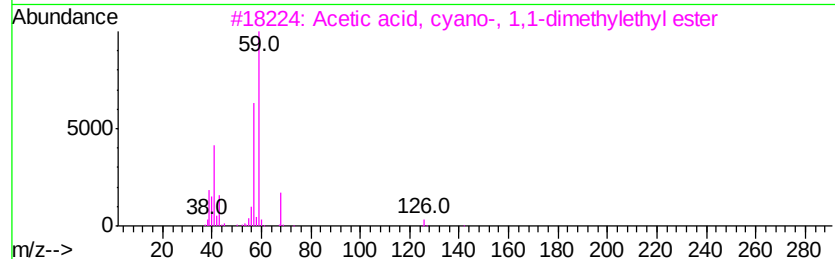
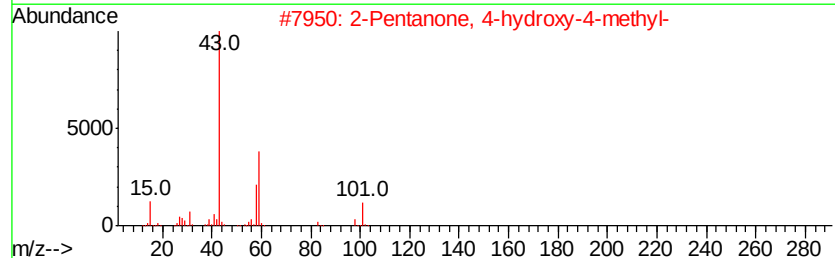
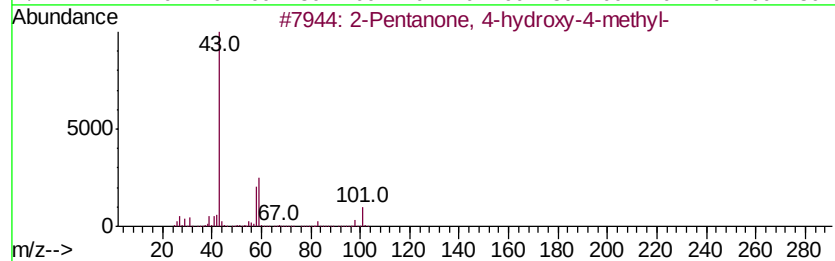
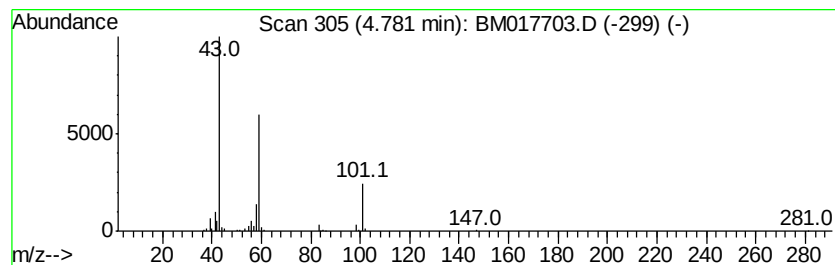
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM11918MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.78	5.37 ng/ul	318940	1,4-Dichlorobenzene-d4	7.60

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	35
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17



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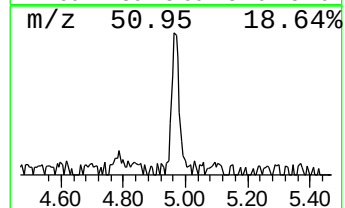
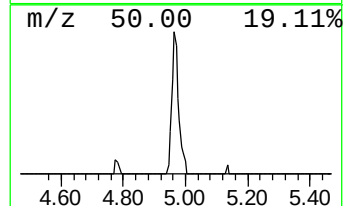
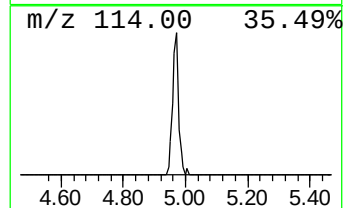
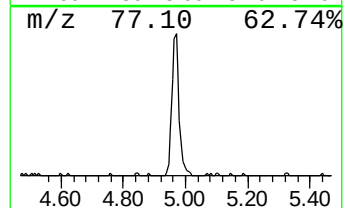
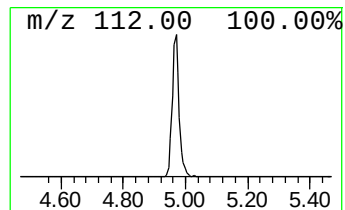
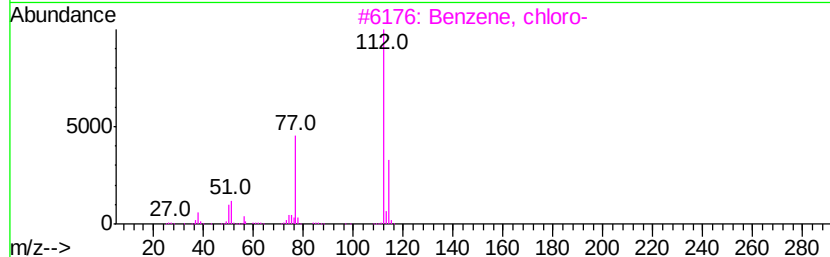
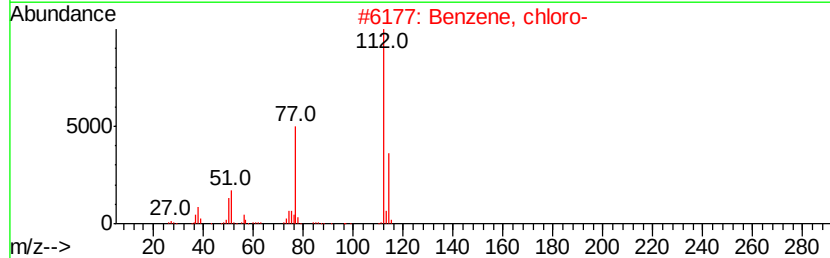
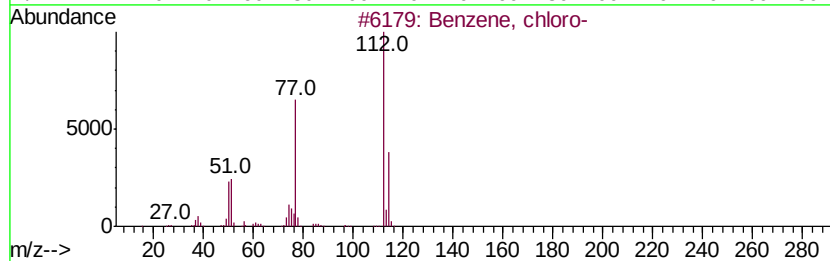
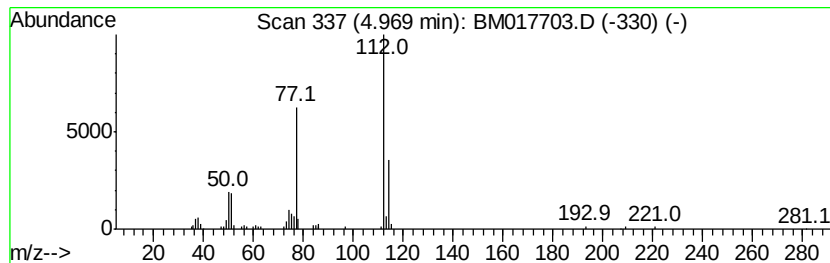
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Benzene, chloro- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.97	2.58 ng/ul	153081	1,4-Dichlorobenzene-d4	7.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, chloro-	112	C6H5Cl	000108-90-7	95
2		Benzene, chloro-	112	C6H5Cl	000108-90-7	94
3		Benzene, chloro-	112	C6H5Cl	000108-90-7	94
4		Benzene, chloro-	112	C6H5Cl	000108-90-7	87
5		2-Chloro-3-nitropyridine	158	C5H3ClN2O2	005470-18-8	42



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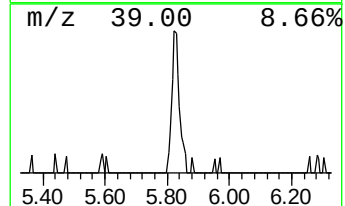
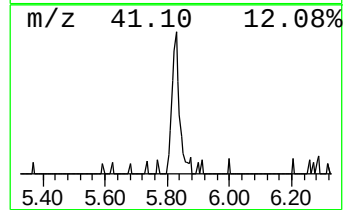
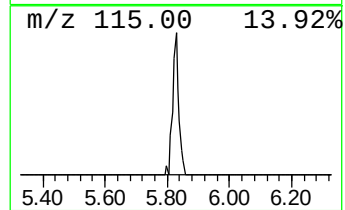
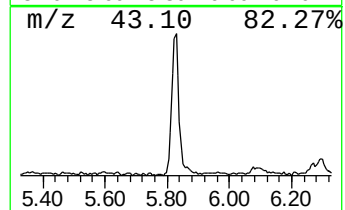
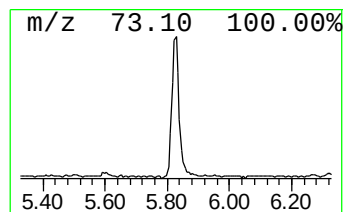
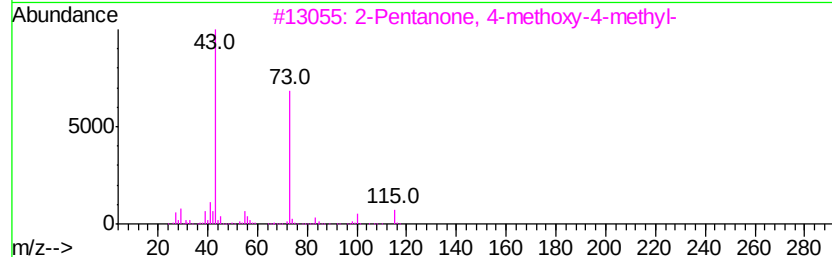
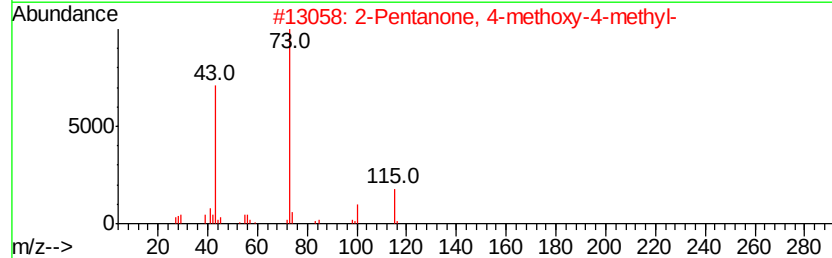
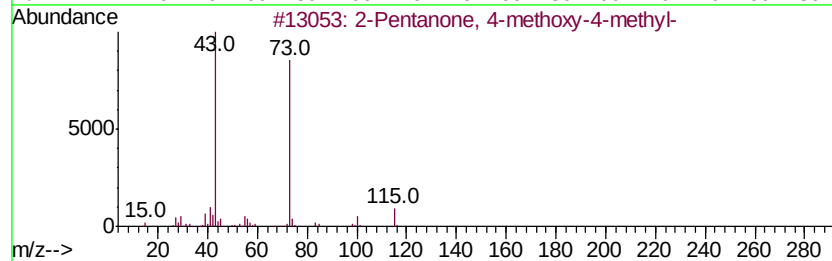
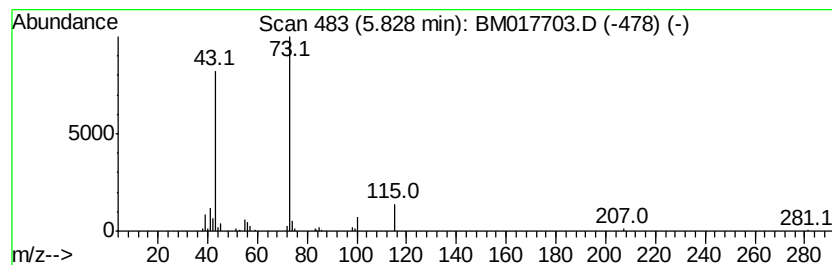
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Peak Number 3 2-Pentanone, 4-methoxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.83	2.49 ng/ul	147898	1,4-Dichlorobenzene-d4	7.60

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	78
2		2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	78
3		2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	50
4		2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	38
5		N-.alpha.-Acetylglycinamide	116	C4H8N2O2	002620-63-5	28



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Data File : BM017703.D
Acq On : 20 Nov 2018 14:11
Operator : MJ/SJ
Sample : J6001-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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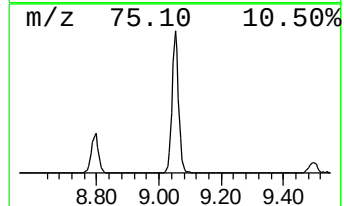
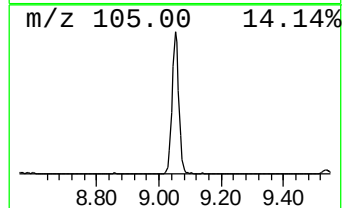
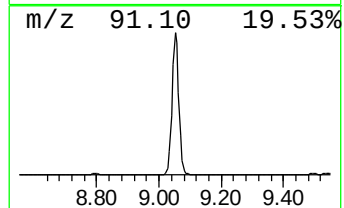
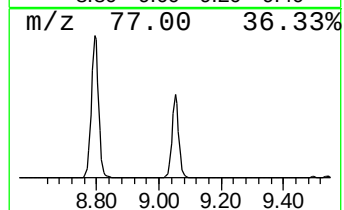
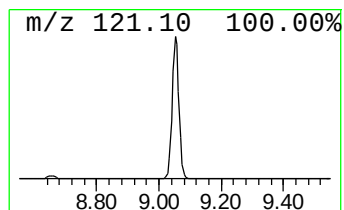
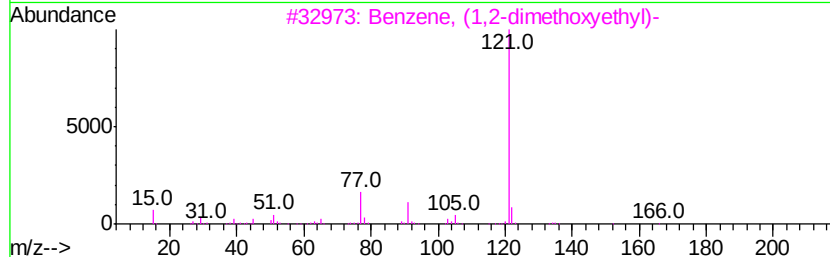
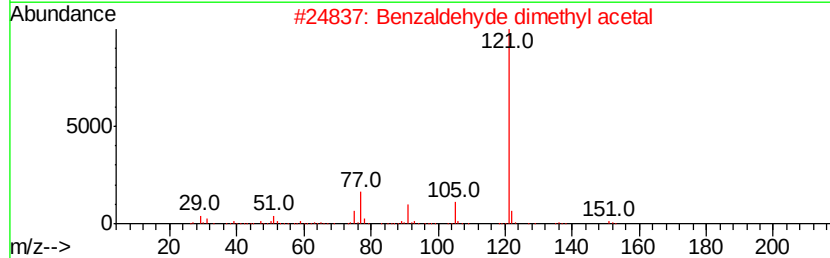
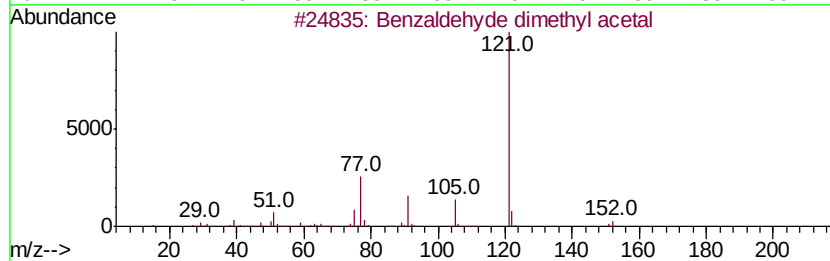
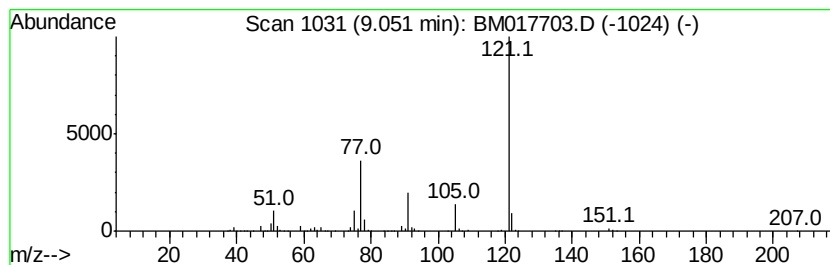
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzaldehyde dimethyl acetal Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.05	38.92 ng/ul	3640570	Naphthalene-d8	10.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzaldehyde dimethyl acetal	152	C9H12O2	001125-88-8	90
2		Benzaldehyde dimethyl acetal	152	C9H12O2	001125-88-8	80
3		Benzene, (1,2-dimethoxyethyl)-	166	C10H14O2	004013-37-0	72
4		Benzeneacetic acid, .alpha.-meth...	180	C10H12O3	056143-21-6	72
5		Benzeneacetic acid, .alpha.-meth...	166	C9H10O3	007021-09-2	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112018\
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Misc :
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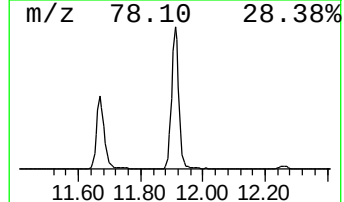
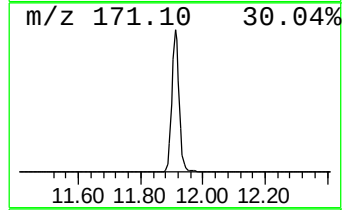
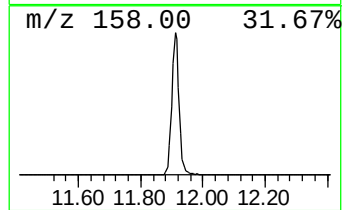
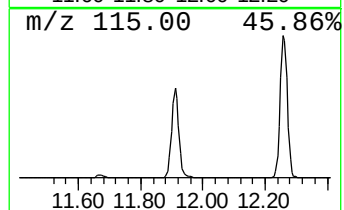
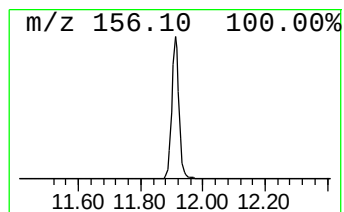
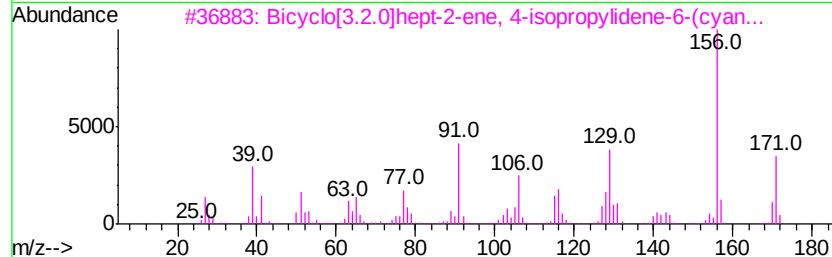
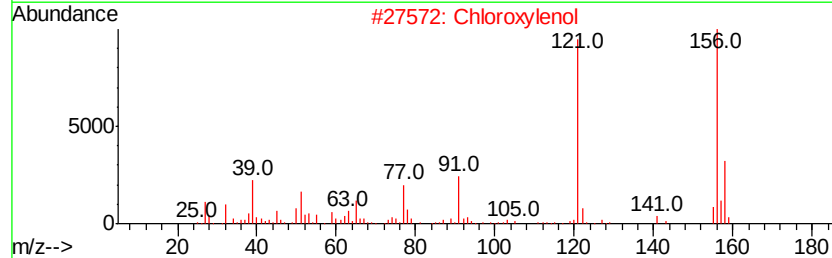
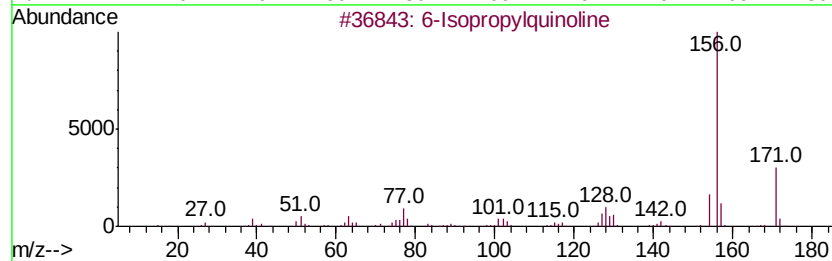
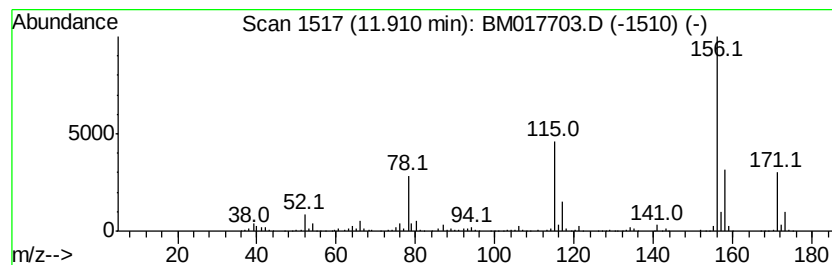
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.91	35.11 ng/ul	3284260	Naphthalene-d8	10.37

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Isopropylquinoline		171	C12H13N	000135-79-5	32
2	Chloroxyleneol		156	C8H9ClO	000088-04-0	32
3	Bicyclo[3.2.0]hept-2-ene, 4-isopropylidene-6-(cyan...		171	C12H13N	1000217-15-6	32
4	Benzene, 2,5-cyclohexadien-1-yl-		156	C12H12	004794-05-2	28
5	1,3-Diaza-5,6-dedihydrocyclohexa...		171	C6H9N3OS	021263-85-4	23



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112018\
Data File : BM017703.D
Acq On : 20 Nov 2018 14:11
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Sample : J6001-01
Misc :
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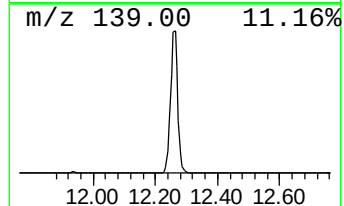
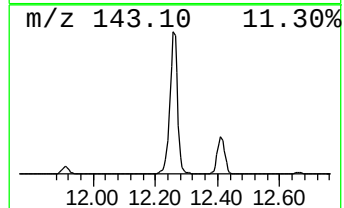
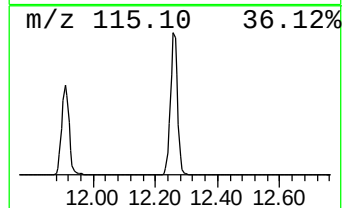
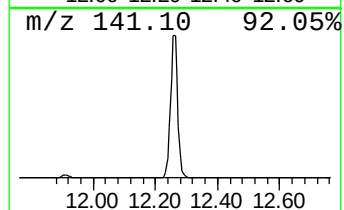
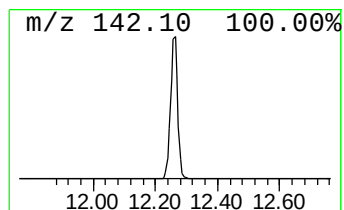
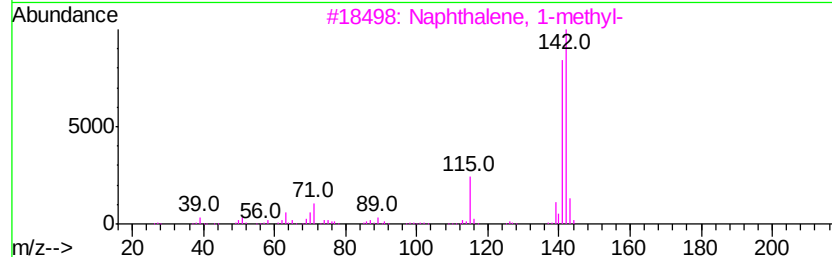
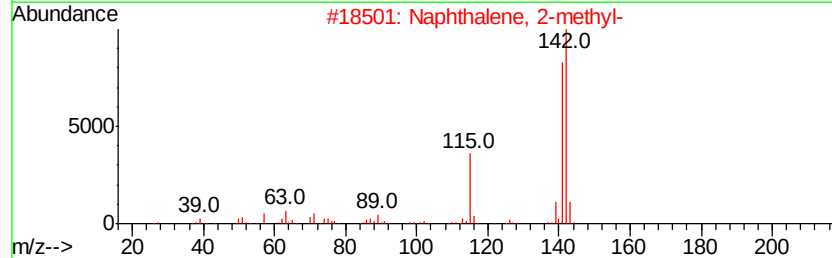
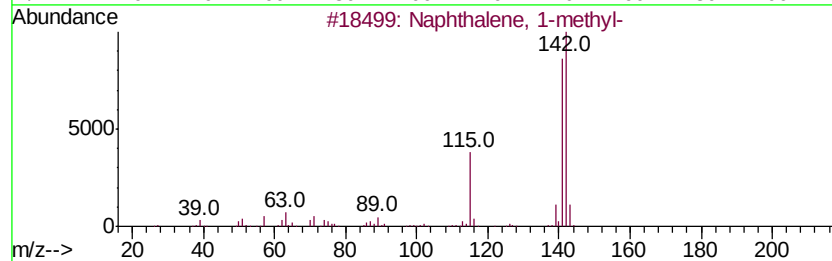
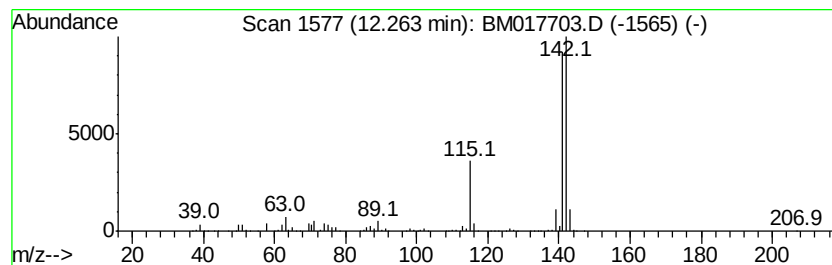
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.26	56.08 ng/ul	5245860	Naphthalene-d8	10.37

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
4		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91
5		Benzocycloheptatriene	142	C11H10	000264-09-5	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112018\
Data File : BM017703.D
Acq On : 20 Nov 2018 14:11
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Sample : J6001-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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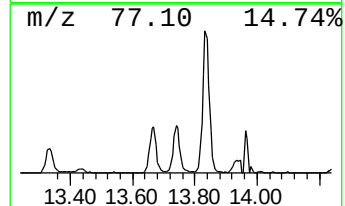
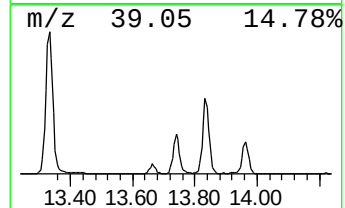
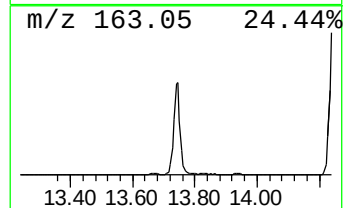
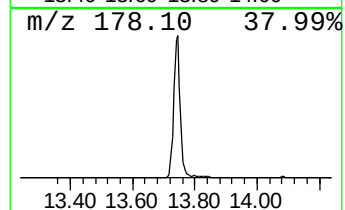
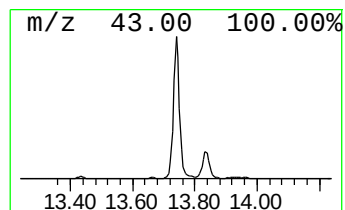
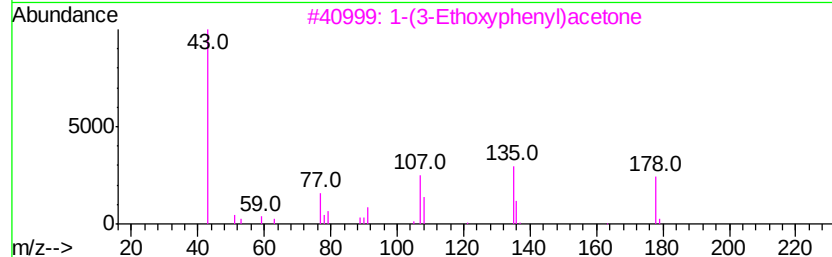
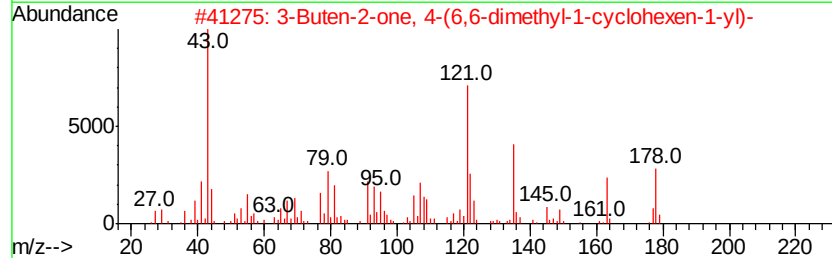
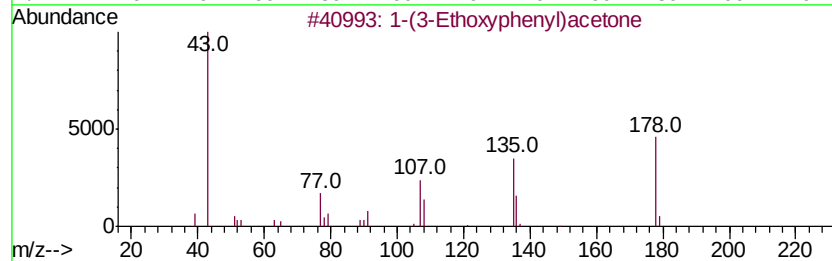
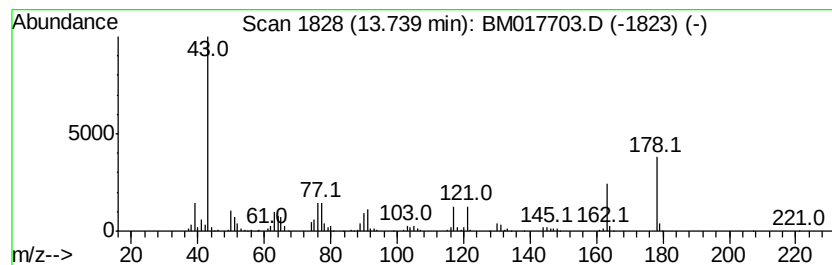
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 unknown-02 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.74	8.34 ng/ul	1064180	Acenaphthene-d10	14.24

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-(3-Ethoxyphenyl)acetone	178	C11H14O2	023037-47-0	30
2		3-Buten-2-one, 4-(6,6-dimethyl-1...	178	C12H18O	065133-79-1	25
3		1-(3-Ethoxyphenyl)acetone	178	C11H14O2	023037-47-0	16
4		2,6-Diacetylpyridine	163	C9H9NO2	001129-30-2	10
5		Carbonodithioic acid, O,S-bis(1-...	178	C7H14OS2	019615-06-6	9



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112018\
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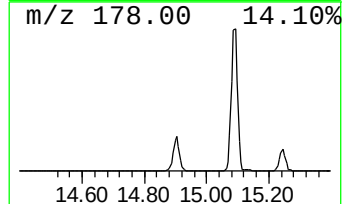
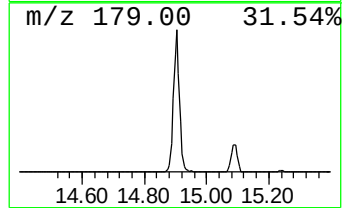
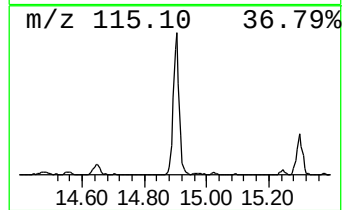
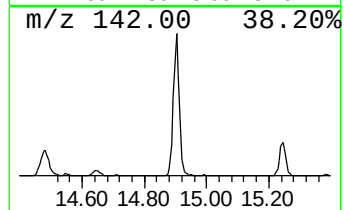
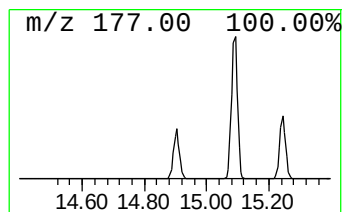
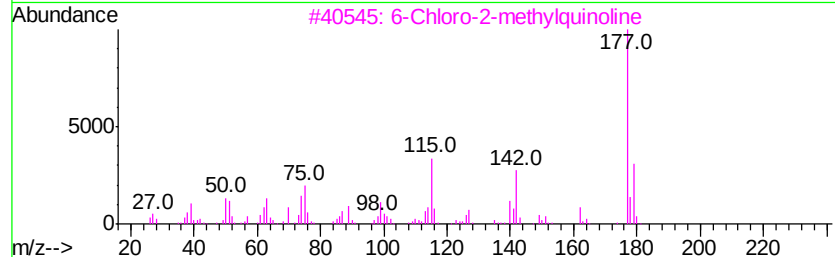
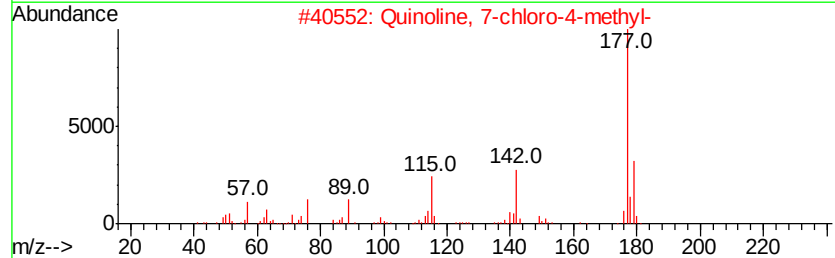
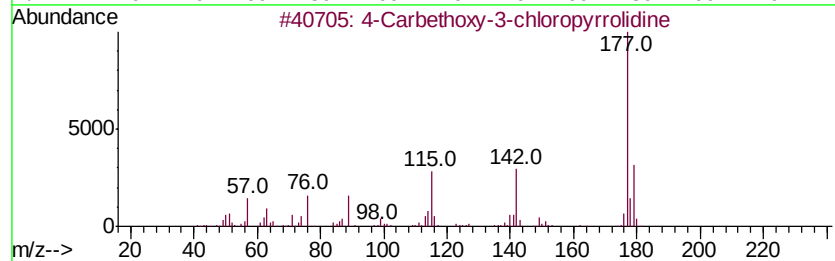
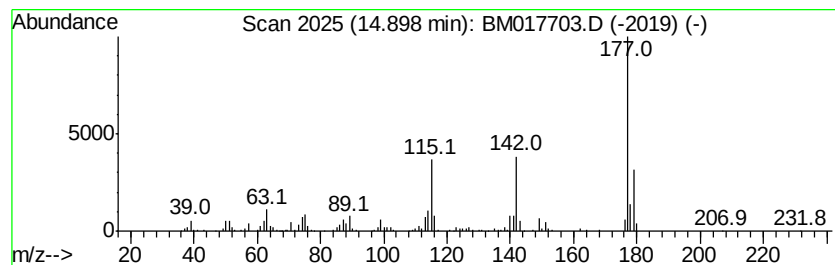
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 4-Carboxy-3-chloropyrrol... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.90	6.10 ng/ul	777704	Acenaphthene-d10	14.24
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		4-Carboxy-3-chloropyrrolidine	177 C7H12ClN02	1000255-97-9 97
2		Quinoline, 7-chloro-4-methyl-	177 C10H8ClN	040941-53-5 91
3		6-Chloro-2-methylquinoline	177 C10H8ClN	000092-46-6 91
4		Quinoline, 2-chloro-4-methyl-	177 C10H8ClN	000634-47-9 90
5		Quinoline, 7-chloro-2-methyl-	177 C10H8ClN	004965-33-7 87



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112018\
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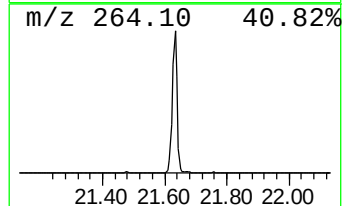
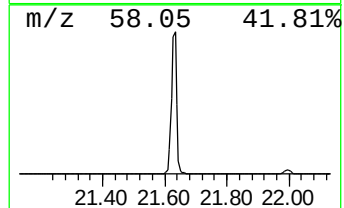
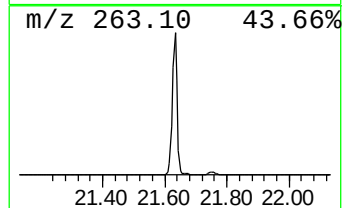
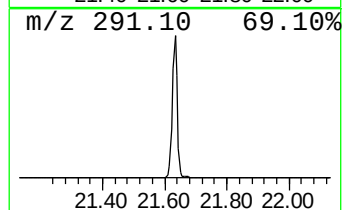
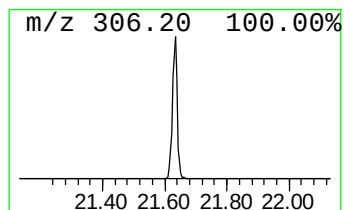
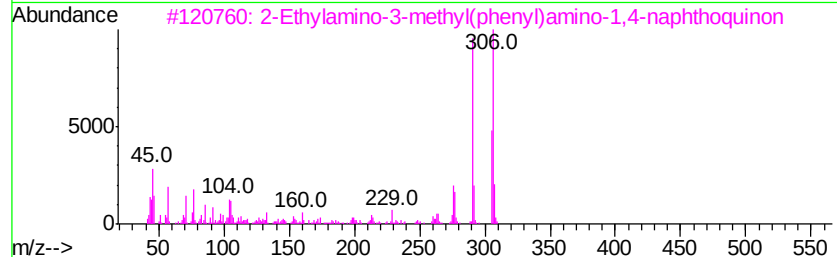
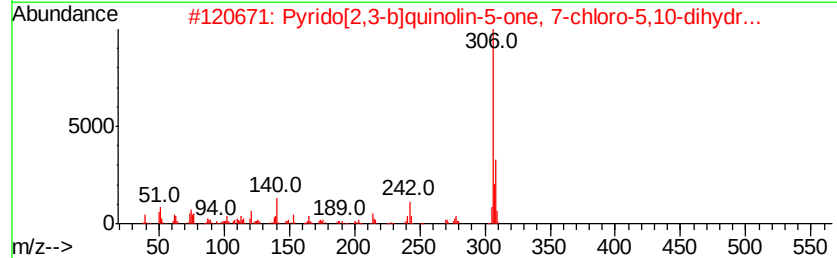
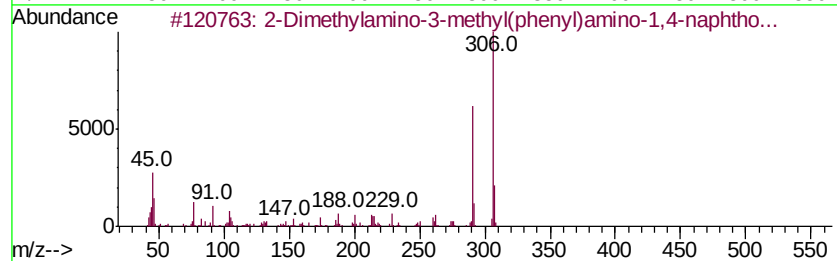
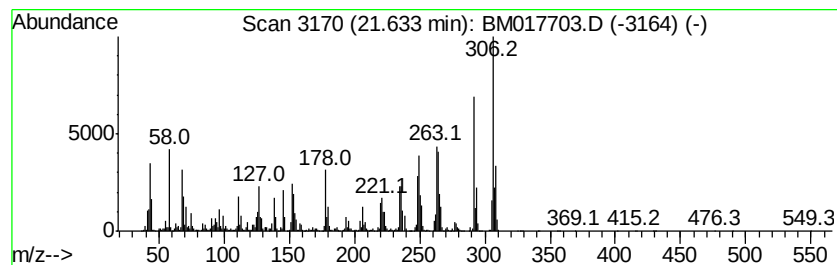
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 unknown-03 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
21.63	15.16 ng/ul	9633740	Chrysene-d12	21.20		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Dimethylamino-3-methyl(phenyl)...	306	C19H18N2O2	134614-10-1	38	
2	Pyrido[2,3-b]quinolin-5-one, 7-c...	306	C18H11ClN2O	069750-07-8	38	
3	2-Ethylamino-3-methyl(phenyl)ami...	306	C19H18N2O2	134614-16-7	35	
4	Azuleno[6,5-b]furan-2(3H)-one, 6...	366	C19H26O7	051292-55-8	27	
5	3,4,5-Trimethoxy-2',4'-diaminodi...	306	C15H18N2O3S	024891-45-0	22	



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112018\
Data File : BM017703.D
Acq On : 20 Nov 2018 14:11
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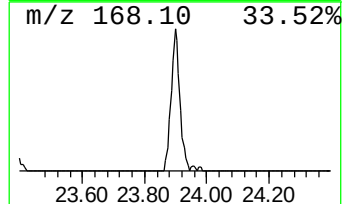
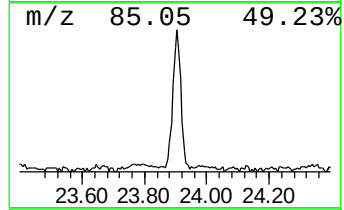
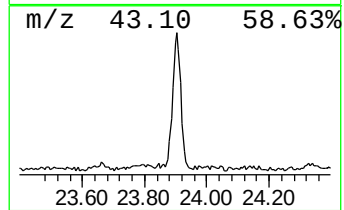
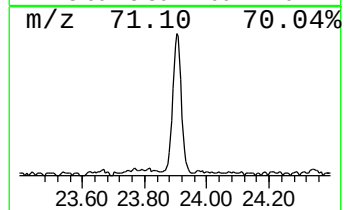
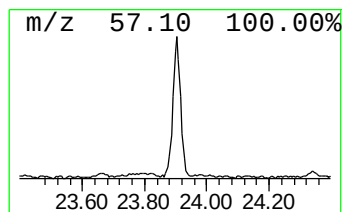
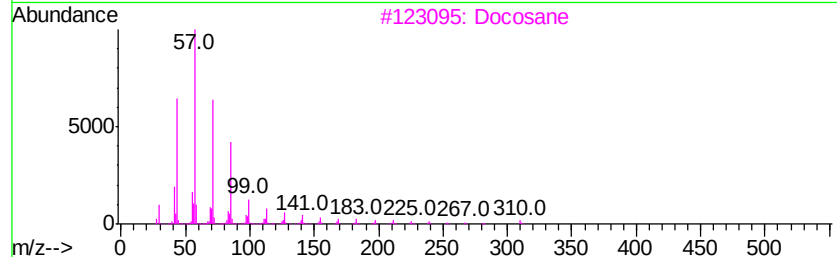
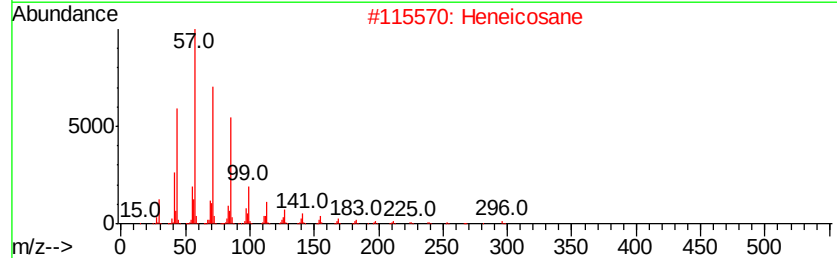
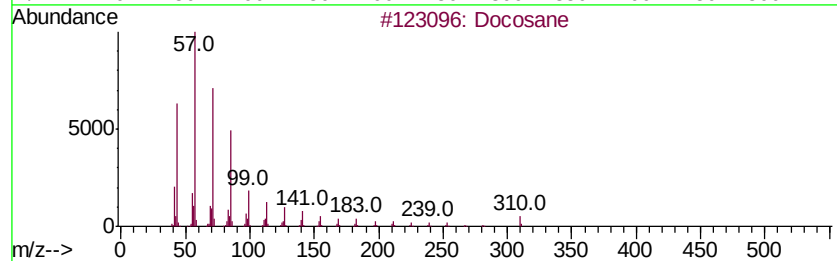
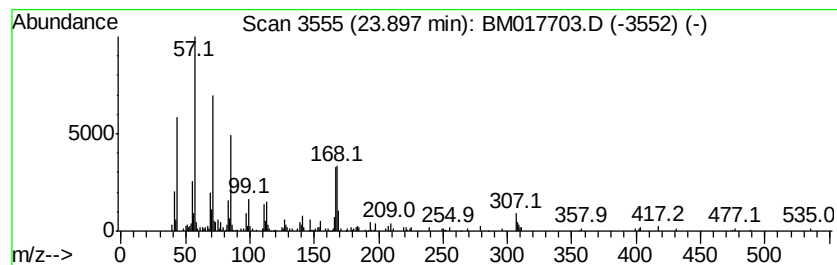
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111918MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.90	2.41 ng/ul	426316	Perylene-d12	23.39

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Docosane	310	C22H46	000629-97-0	91
2		Heneicosane	296	C21H44	000629-94-7	89
3		Docosane	310	C22H46	000629-97-0	70
4		Eicosane	282	C20H42	000112-95-8	70
5		Decane, 3,6-dimethyl-	170	C12H26	017312-53-7	55



Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM112018\
Data File : BM017703.D
Acq On : 20 Nov 2018 14:11
Operator : MJ/SJ
Sample : J6001-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A40T1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111918MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.78	5.4	ng/ul	318940	1	7.60	1188370	20.0
Benzene, chloro-	4.97	2.6	ng/ul	153081	1	7.60	1188370	20.0
2-Pentanone, 4-me...	5.83	2.5	ng/ul	147898	1	7.60	1188370	20.0
Benzaldehyde dime...	9.05	38.9	ng/ul	3640570	2	10.37	1870810	20.0
unknown-01	11.91	35.1	ng/ul	3284260	2	10.37	1870810	20.0
Naphthalene, 1-me...	12.26	56.1	ng/ul	5245860	2	10.37	1870810	20.0
unknown-02	13.74	8.3	ng/ul	1064180	3	14.24	2550880	20.0
4-Carbethoxy-3-ch...	14.90	6.1	ng/ul	777704	3	14.24	2550880	20.0
unknown-03	21.63	15.2	ng/ul	9633740	5	21.20	12708100	20.0
(DEL) Alkane: Str...	23.90	2.4	ng/ul	426316	6	23.39	3545180	20.0