

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111319\
 Data File : BN008567.D
 Acq On : 14 Nov 2019 21:31
 Operator : JU
 Sample : K5739-03
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0CR1

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_N\METHODS\SOM-EPA-BN111319MA.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	3.222	36	40	49	rVB	80936	121478	1.38%	0.106%
2	3.840	142	145	150	rVB2	21406	33166	0.38%	0.029%
3	5.011	337	344	350	rBV	134860	207423	2.36%	0.182%
4	6.816	644	651	662	rBV3	359222	722944	8.23%	0.633%
5	6.975	672	678	691	rBV	935615	1490955	16.98%	1.305%
6	7.175	703	712	721	rBV	1163330	1831702	20.86%	1.603%
7	7.269	724	728	735	rVV2	31736	55074	0.63%	0.048%
8	7.528	765	772	778	rVB2	37980	59696	0.68%	0.052%
9	7.634	783	790	794	rVV	1205574	1901046	21.65%	1.664%
10	7.669	794	796	802	rVV	275144	356313	4.06%	0.312%
11	7.981	842	849	856	rVB	1224294	1953354	22.25%	1.709%
12	8.334	902	909	920	rBV	938535	1505158	17.14%	1.317%
13	8.452	922	929	932	rVB6	15300	25013	0.28%	0.022%
14	8.775	977	984	986	rVV	1277231	2096791	23.88%	1.835%
15	8.816	986	991	1002	rVV	5371316	8779008	100.00%	7.682%
16	9.228	1054	1061	1064	rBV2	48218	78537	0.89%	0.069%
17	9.263	1064	1067	1073	rVB3	49679	74600	0.85%	0.065%
18	9.487	1098	1105	1119	rVV	1139211	1950761	22.22%	1.707%
19	10.022	1189	1196	1205	rBV	1664688	2739999	31.21%	2.398%
20	10.181	1217	1223	1229	rVB9	25543	49950	0.57%	0.044%
21	10.263	1231	1237	1245	rBV	593803	1008076	11.48%	0.882%
22	10.398	1252	1260	1267	rBV	1566559	2644049	30.12%	2.314%
23	10.528	1274	1282	1292	rVB	113652	207982	2.37%	0.182%
24	10.781	1316	1325	1331	rBV	561291	915703	10.43%	0.801%
25	10.987	1349	1360	1371	rBV	4498375	7355712	83.79%	6.437%
26	11.198	1388	1396	1404	rBV	1110618	1821275	20.75%	1.594%
27	11.687	1473	1479	1486	rBV2	166863	280154	3.19%	0.245%
28	11.751	1486	1490	1494	rVB2	34127	50070	0.57%	0.044%
29	11.993	1524	1531	1535	rBV2	31042	53744	0.61%	0.047%
30	12.375	1585	1596	1603	rBV	231477	399287	4.55%	0.349%
31	12.434	1603	1606	1613	rVB3	31059	44817	0.51%	0.039%
32	12.669	1633	1646	1651	rVB6	77579	160427	1.83%	0.140%
33	12.757	1656	1661	1666	rBV8	23501	40203	0.46%	0.035%
34	12.822	1666	1672	1679	rVV	192540	283023	3.22%	0.248%

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Integration Parameters: LSCINT.P

Integrator: RTE
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 Sampling : 1
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 Title : SVOA CALIBRATION

35	12.998	1697	1702	1705	rBV4	16350	25817	0.29%	0.023%
36	13.051	1705	1711	1719	rVB2	398858	705204	8.03%	0.617%
37	13.151	1724	1728	1732	rVV4	79821	126283	1.44%	0.111%
38	13.204	1732	1737	1742	rVV3	98798	176678	2.01%	0.155%
39	13.269	1742	1748	1753	rVV2	138035	255116	2.91%	0.223%
40	13.322	1753	1757	1763	rVB4	19761	29618	0.34%	0.026%
41	13.639	1804	1811	1813	rBV2	382508	547821	6.24%	0.479%
42	13.675	1813	1817	1822	rVV	3332466	4535744	51.67%	3.969%
43	13.722	1822	1825	1830	rVV	319793	445298	5.07%	0.390%
44	13.763	1830	1832	1839	rVB7	16376	26036	0.30%	0.023%
45	13.904	1851	1856	1857	rBV4	53661	74070	0.84%	0.065%
46	13.945	1857	1863	1870	rVV	3453626	5055121	57.58%	4.423%
47	14.034	1872	1878	1885	rVB6	61211	116315	1.32%	0.102%
48	14.257	1909	1916	1926	rVB	2489578	3649615	41.57%	3.194%
49	14.451	1941	1949	1958	rVV	413615	670808	7.64%	0.587%
50	14.534	1959	1963	1968	rVV2	48977	77789	0.89%	0.068%
51	14.739	1994	1998	2004	rVV2	114678	167720	1.91%	0.147%
52	15.016	2042	2045	2050	rBV	53237	71917	0.82%	0.063%
53	15.151	2063	2068	2072	rBV4	16864	25536	0.29%	0.022%
54	15.251	2079	2085	2096	rBV	4197576	6184723	70.45%	5.412%
55	15.369	2099	2105	2116	rBV	1428906	2011025	22.91%	1.760%
56	15.492	2122	2126	2131	rBV2	60133	81401	0.93%	0.071%
57	15.828	2179	2183	2187	rVV6	57900	80196	0.91%	0.070%
58	15.916	2192	2198	2201	rVV8	19381	43092	0.49%	0.038%
59	15.963	2203	2206	2209	rVV5	21819	31765	0.36%	0.028%
60	15.998	2209	2212	2215	rVV5	34651	50344	0.57%	0.044%
61	16.033	2215	2218	2225	rVB4	56996	87520	1.00%	0.077%
62	16.122	2230	2233	2236	rVV5	21268	26197	0.30%	0.023%
63	16.228	2246	2251	2256	rVB2	201239	278024	3.17%	0.243%
64	16.275	2256	2259	2264	rVB3	25102	34667	0.39%	0.030%
65	16.363	2270	2274	2276	rBV5	19382	27491	0.31%	0.024%
66	16.398	2276	2280	2284	rVB6	66406	83556	0.95%	0.073%
67	16.563	2303	2308	2315	rVB5	100154	171332	1.95%	0.150%
68	16.810	2342	2350	2357	rVB3	62092	122568	1.40%	0.107%
69	16.998	2376	2382	2389	rBV	2941594	4118148	46.91%	3.604%
70	17.098	2392	2399	2408	rBV	5051800	6890516	78.49%	6.030%
71	17.192	2411	2415	2419	rBV6	72280	122353	1.39%	0.107%

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Integration Parameters: LSCINT.P

Integrator: RTE
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72	17.233	2419	2422	2431	rVB6	92451	187074	2.13%	0.164%
73	17.363	2441	2444	2448	rVB2	35697	45165	0.51%	0.040%
74	17.598	2478	2484	2485	rVV6	27191	51438	0.59%	0.045%
75	17.616	2485	2487	2492	rVB4	63293	82669	0.94%	0.072%
76	17.710	2497	2503	2509	rVV9	94001	152894	1.74%	0.134%
77	17.927	2536	2540	2543	rBV3	96528	128952	1.47%	0.113%
78	17.969	2543	2547	2556	rVB6	87118	188518	2.15%	0.165%
79	18.080	2562	2566	2569	rBV6	26704	34265	0.39%	0.030%
80	18.375	2606	2616	2624	rBV	3651659	4928813	56.14%	4.313%
81	18.451	2626	2629	2633	rVV5	57693	77375	0.88%	0.068%
82	18.504	2633	2638	2645	rVB3	170579	264002	3.01%	0.231%
83	18.669	2662	2666	2669	rBV5	83239	137242	1.56%	0.120%
84	18.704	2669	2672	2676	rVV3	194301	262896	2.99%	0.230%
85	18.751	2676	2680	2685	rVB7	56099	90524	1.03%	0.079%
86	18.839	2691	2695	2699	rBV7	41214	56203	0.64%	0.049%
87	19.027	2724	2727	2733	rVB	65268	94737	1.08%	0.083%
88	19.280	2764	2770	2774	rBV	57816	112560	1.28%	0.098%
89	19.398	2784	2790	2797	rBV	5669176	7886984	89.84%	6.901%
90	19.916	2874	2878	2881	rBV6	58107	93619	1.07%	0.082%
91	20.945	3049	3053	3058	rBV2	697891	910269	10.37%	0.797%
92	21.198	3091	3096	3101	rBV	3536956	4426085	50.42%	3.873%
93	21.398	3127	3130	3134	rVB	229643	241091	2.75%	0.211%
94	21.833	3200	3204	3210	rBV2	349351	443715	5.05%	0.388%
95	22.286	3277	3281	3287	rBV	489475	632134	7.20%	0.553%
96	22.780	3361	3365	3371	rBV	593511	817476	9.31%	0.715%
97	23.239	3437	3443	3453	rVB	4048479	7237256	82.44%	6.333%
98	23.339	3455	3460	3462	rVV	604549	1044448	11.90%	0.914%
99	23.374	3462	3466	3473	rVB	2374313	4219350	48.06%	3.692%
100	23.968	3562	3567	3573	rVB	517724	906744	10.33%	0.793%

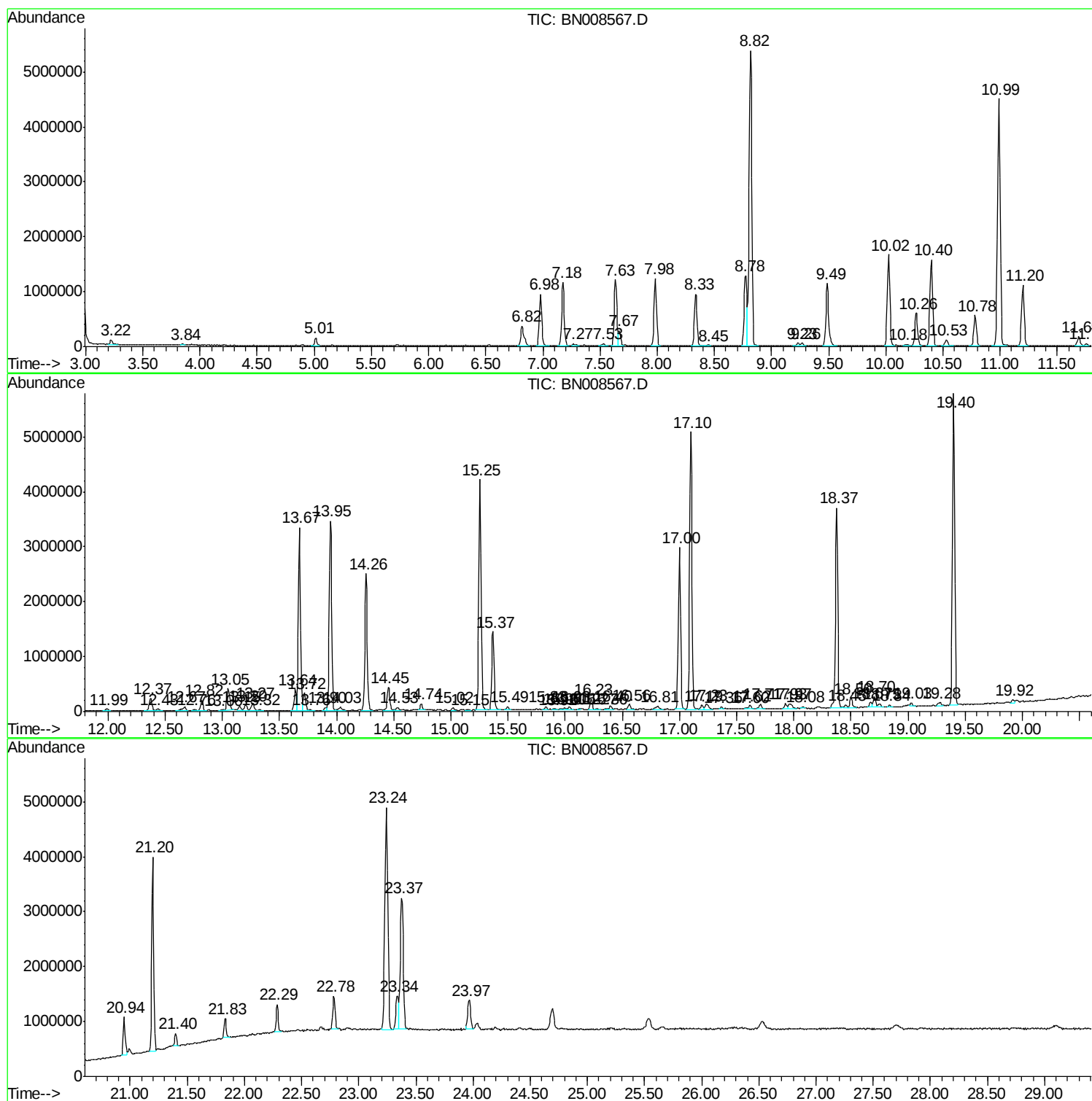
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Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0CR1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN111319MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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Instrument :
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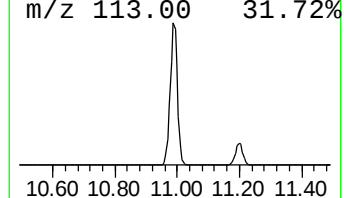
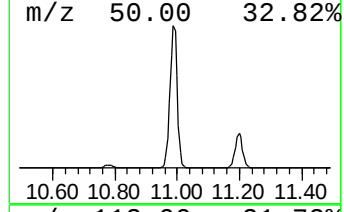
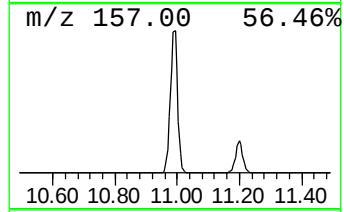
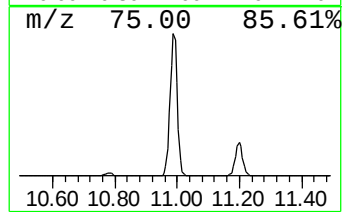
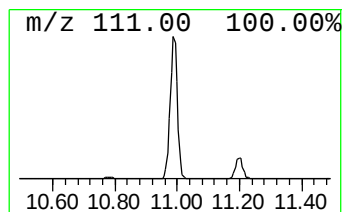
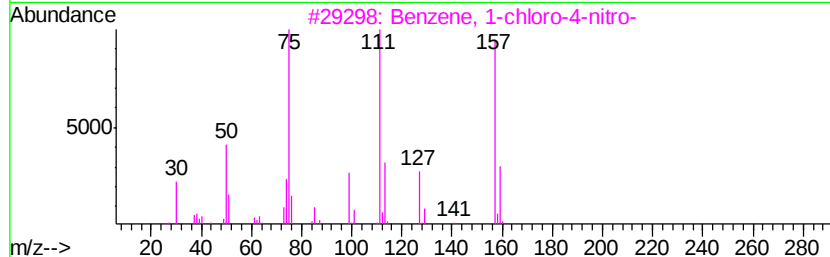
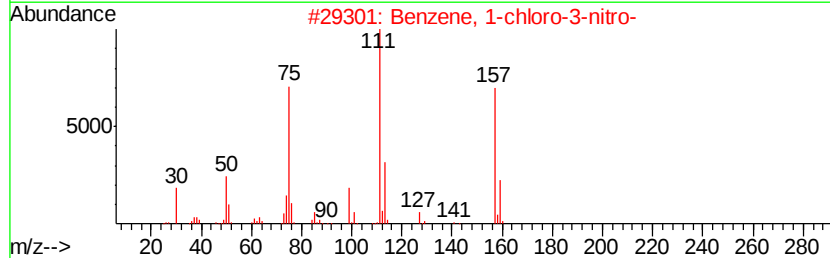
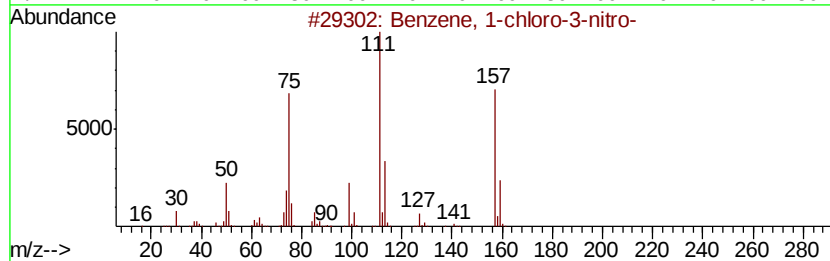
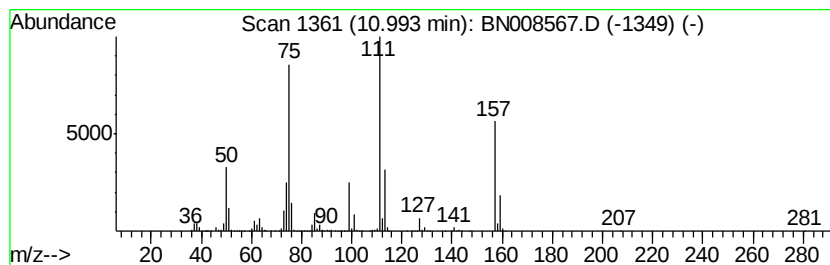
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN111319MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1-chloro-3-nitro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.99	55.64 ng/ul	7355710	Naphthalene-d8	10.40

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	96
2		Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	94
3		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	94
4		Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	93
5		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	87



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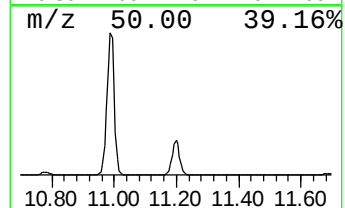
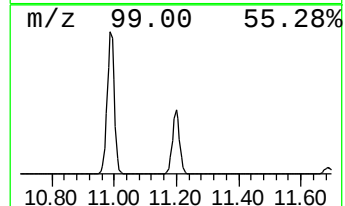
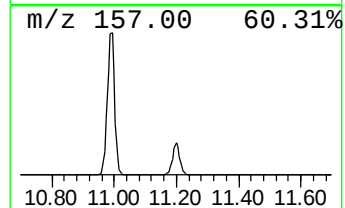
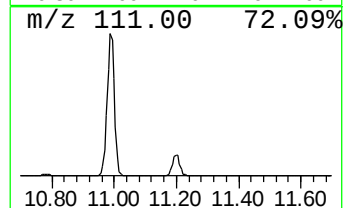
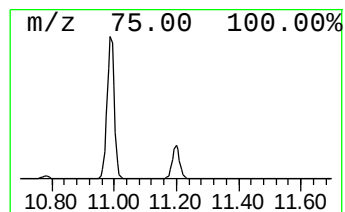
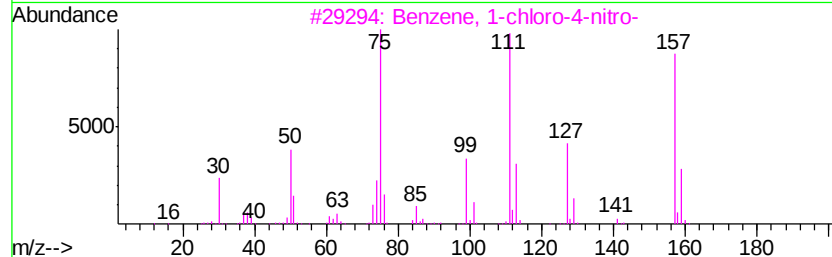
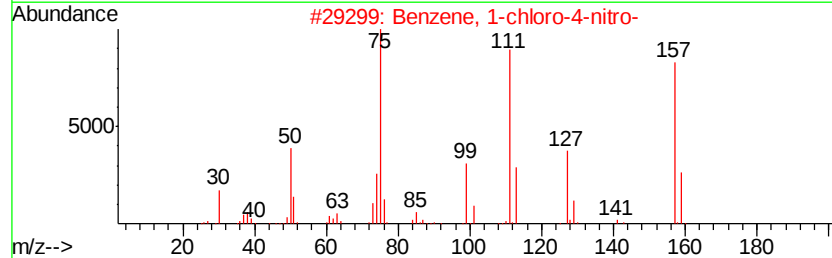
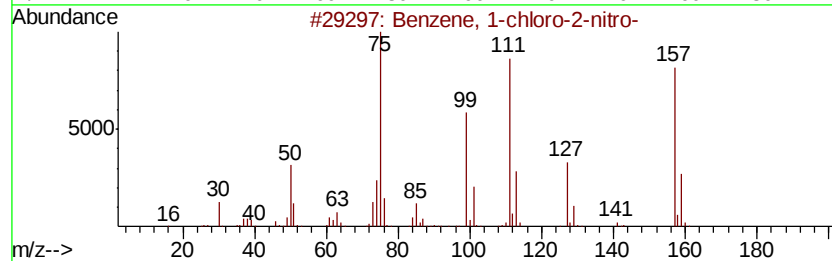
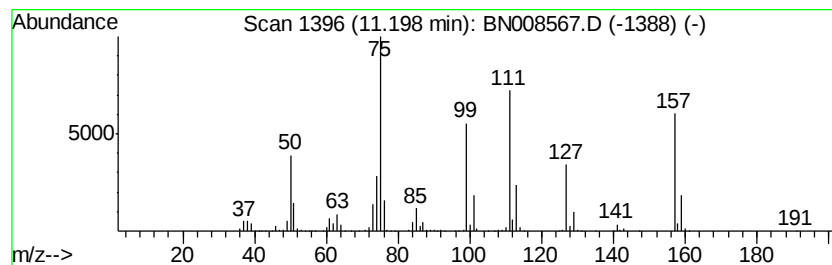
Instrument :
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ClientSampled :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1-chloro-2-nitro- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.20	13.78 ng/ul	1821280	Naphthalene-d8	10.40
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzene, 1-chloro-2-nitro-	157 C6H4ClNO2	000088-73-3	97
2	Benzene, 1-chloro-4-nitro-	157 C6H4ClNO2	000100-00-5	96
3	Benzene, 1-chloro-4-nitro-	157 C6H4ClNO2	000100-00-5	95
4	Benzene, 1-chloro-2-nitro-	157 C6H4ClNO2	000088-73-3	95
5	Benzene, 1-chloro-3-nitro-	157 C6H4ClNO2	000121-73-3	94



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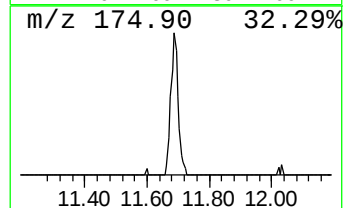
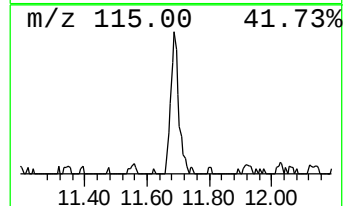
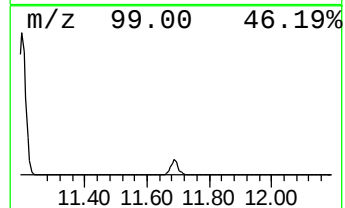
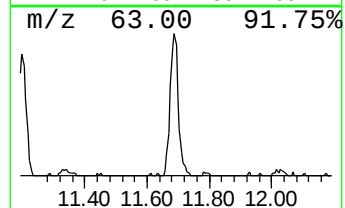
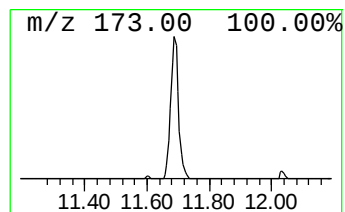
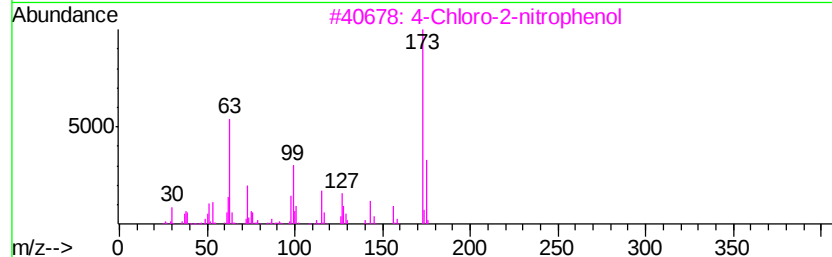
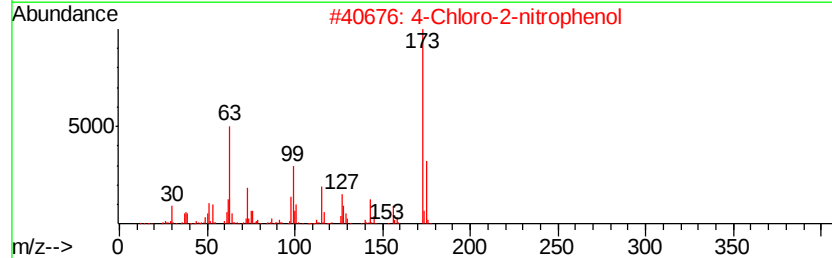
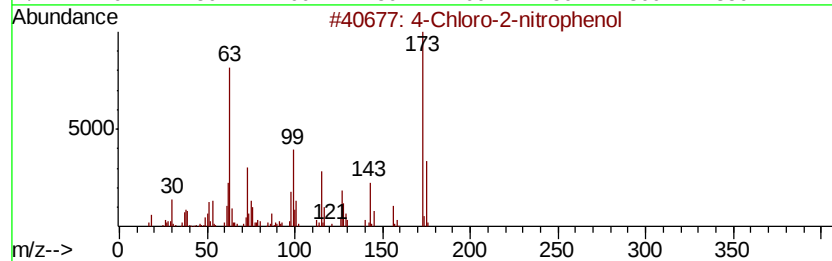
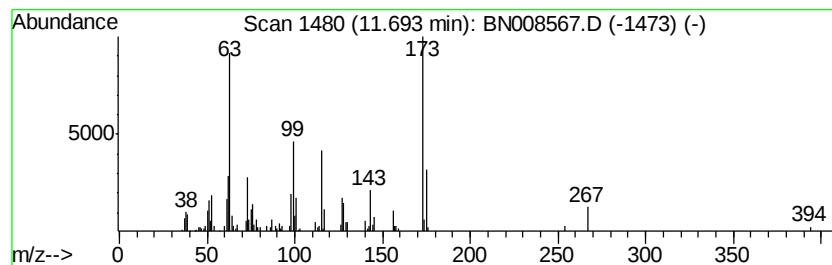
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN111319MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 4-Chloro-2-nitrophenol Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.69	2.12 ng/ul	280154	Naphthalene-d8	10.40

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	99
2		4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	94
3		4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	91
4		Phenol, 5-chloro-2-nitro-	173	C6H4ClNO3	000611-07-4	87
5		4-Chloro-3-nitrophenol	173	C6H4ClNO3	000610-78-6	70



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111319\
Data File : BN008567.D
Acq On : 14 Nov 2019 21:31
Operator : JU
Sample : K5739-03
Misc :
ALS Vial : 44 Sample Multiplier: 1

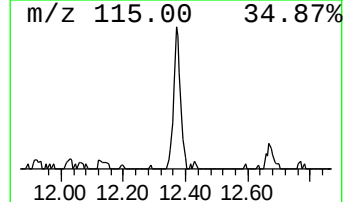
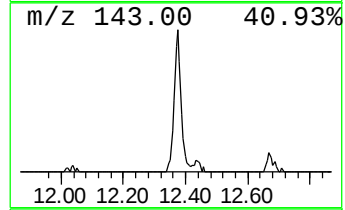
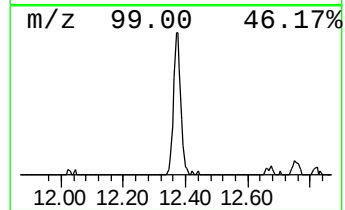
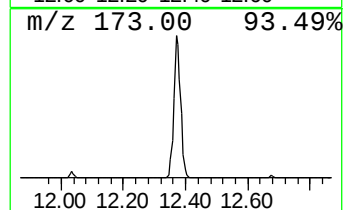
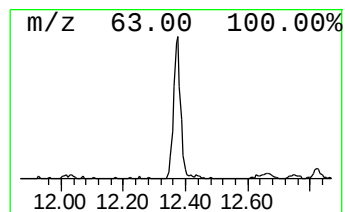
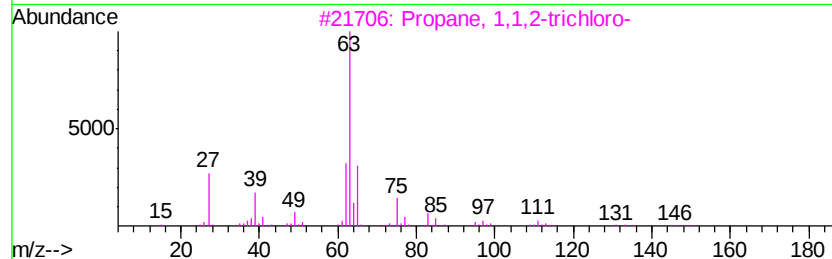
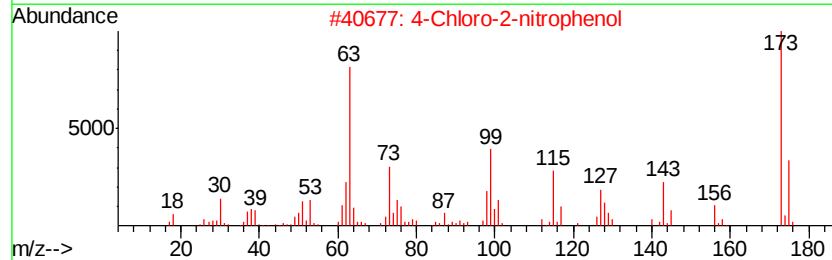
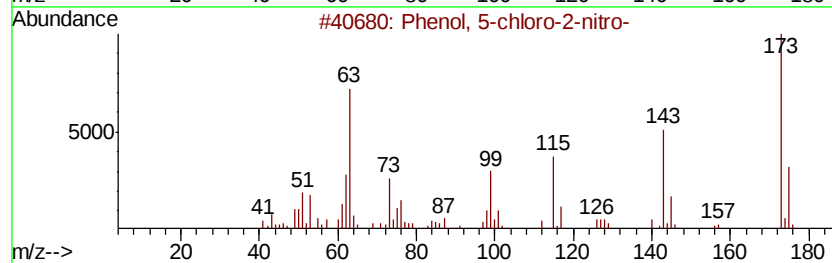
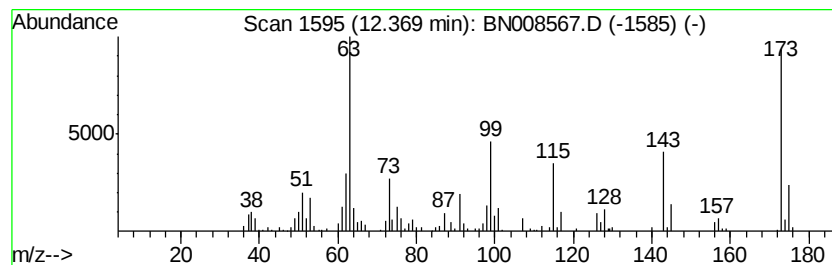
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Phenol, 5-chloro-2-nitro- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.37	2.19 ng/ul	399287	Acenaphthene-d10		14.26	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenol, 5-chloro-2-nitro-	173	C6H4ClNO3	000611-07-4	94
2		4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	55
3		Propane, 1,1,2-trichloro-	146	C3H5Cl3	000598-77-6	47
4		4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	43
5		Propanoic acid, 2-chloro-, ethyl...	136	C5H9ClO2	000535-13-7	40



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111319\
Data File : BN008567.D
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ClientSampleId :
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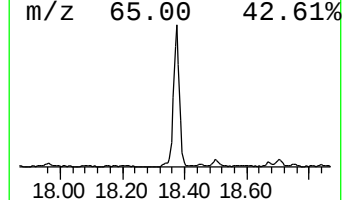
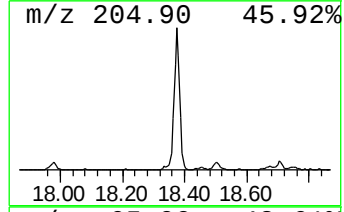
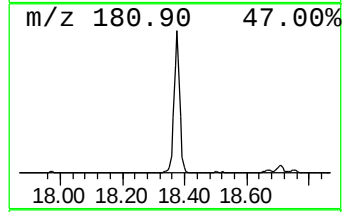
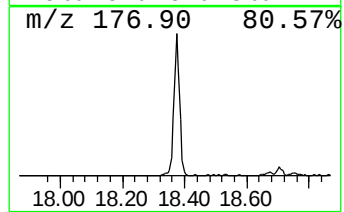
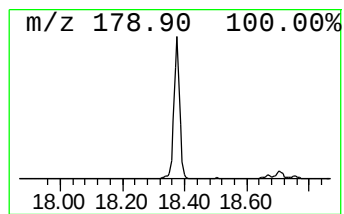
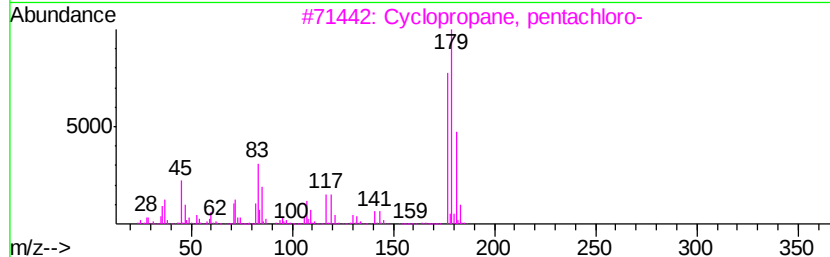
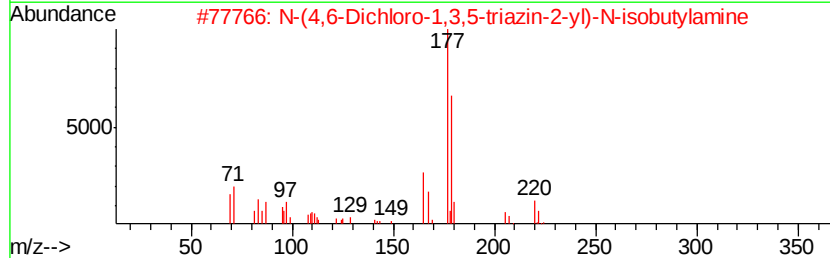
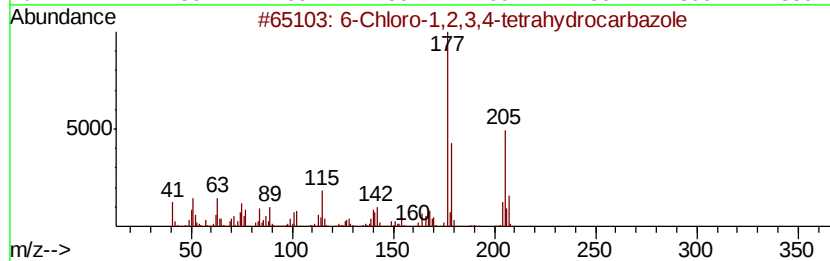
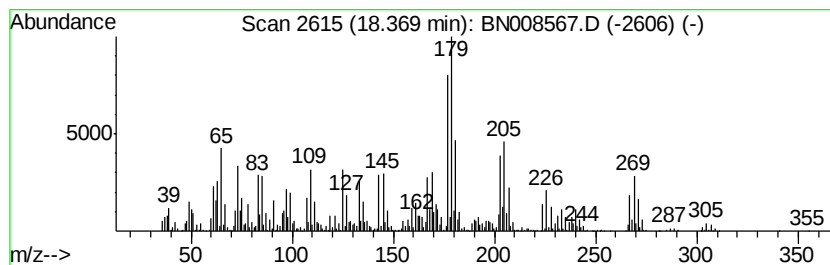
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.37	23.94 ng/ul	4928810	Phenanthrene-d10	17.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	6-Chloro-1,2,3,4-tetrahydrocarba...	205	C12H12ClN	036684-65-8	25
2		N-(4,6-Dichloro-1,3,5-triazin-2-...	220	C7H10Cl2N4	1000326-88-2	12
3		Cyclopropane, pentachloro-	212	C3HCl5	006262-51-7	12
4		1H-Pyrrole, 1-(4-chlorophenyl)-	177	C10H8ClN	005044-38-2	11
5		2,4,6-Trifluoronitrobenzene	177	C6H2F3NO2	000315-14-0	11



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111319\
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Misc :
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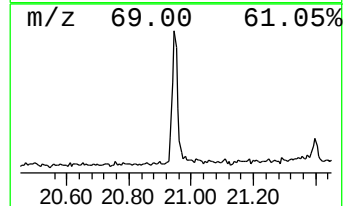
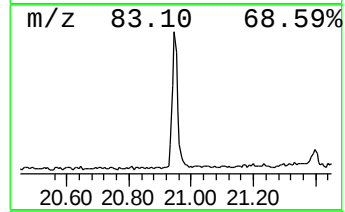
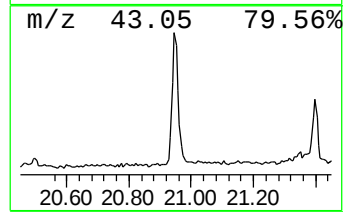
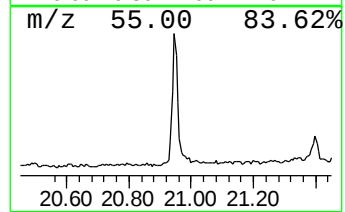
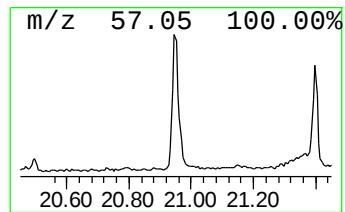
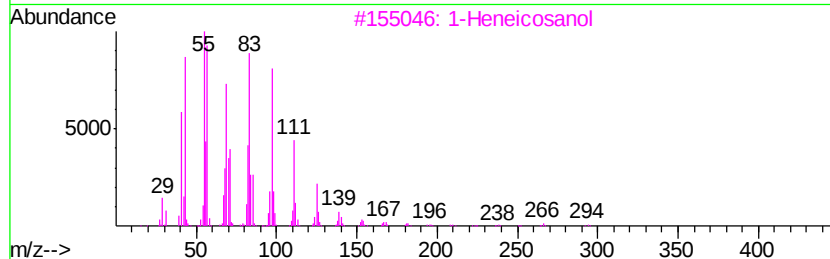
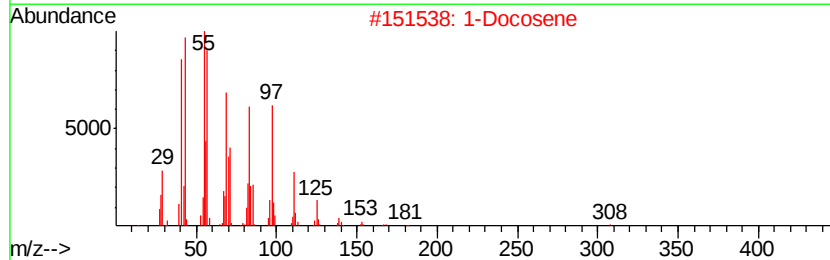
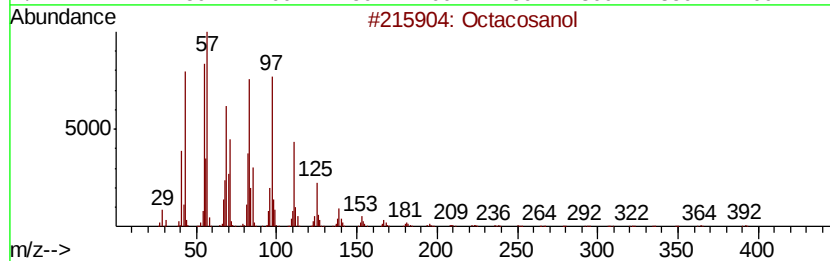
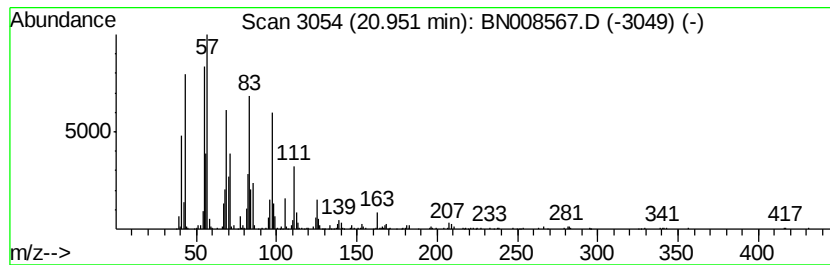
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Octacosanol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.95	4.11 ng/ul	910269	Chrysene-d12	21.20
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Octacosanol	410 C28H58O	000557-61-9	94
2	1-Docosene	308 C22H44	001599-67-3	94
3	1-Heneicosanol	312 C21H44O	015594-90-8	94
4	Pentafluoropropionic acid, hexad...	388 C19H33F5O2	006222-07-7	91
5	1-Tricosene	322 C23H46	018835-32-0	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111319\
Data File : BN008567.D
Acq On : 14 Nov 2019 21:31
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Instrument :
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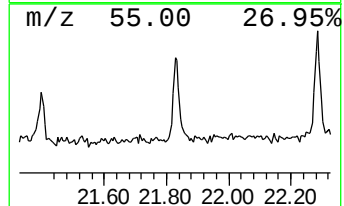
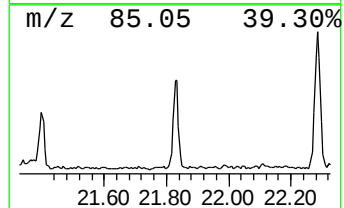
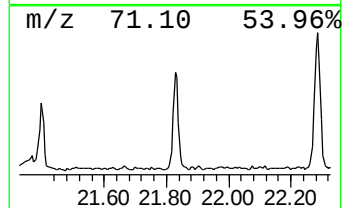
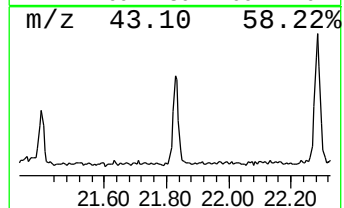
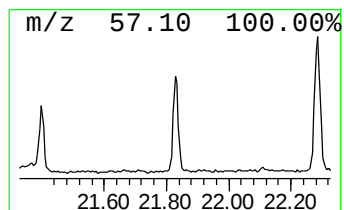
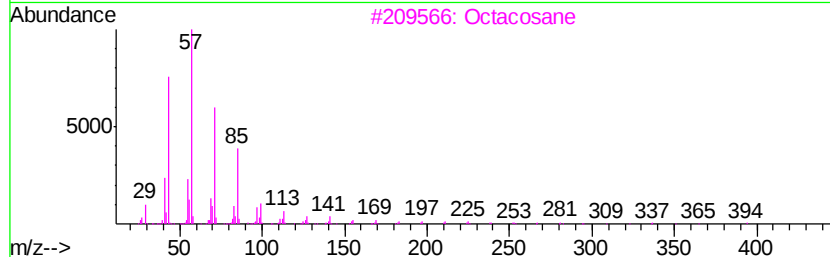
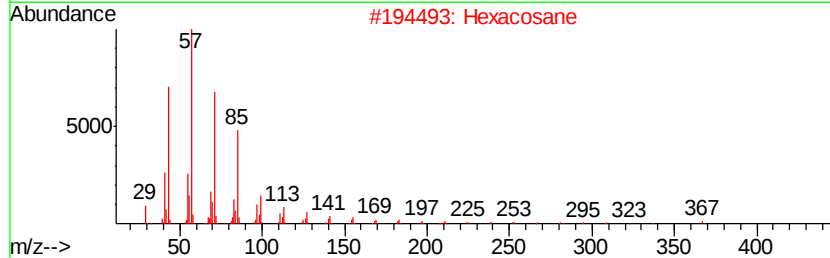
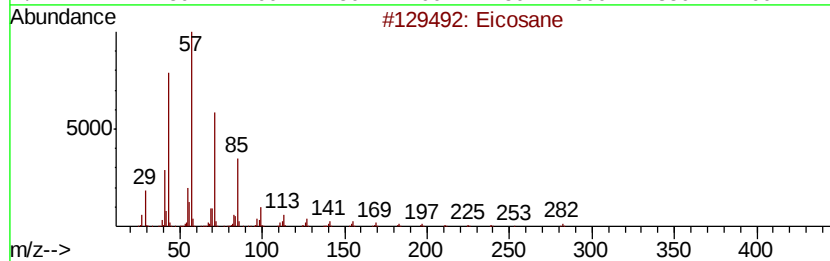
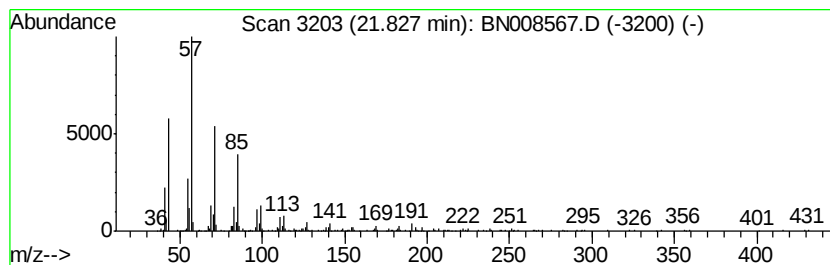
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.83	2.01 ng/ul	443715	Chrysene-d12	21.20

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	93
2		Hexacosane	366	C26H54	000630-01-3	90
3		Octacosane	394	C28H58	000630-02-4	90
4		Tritetracontane	605	C43H88	007098-21-7	90
5		Hentriacontane	437	C31H64	000630-04-6	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111319\
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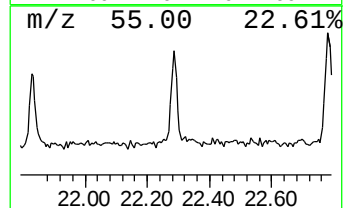
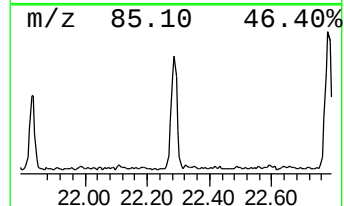
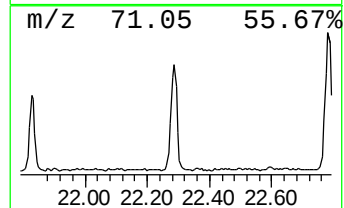
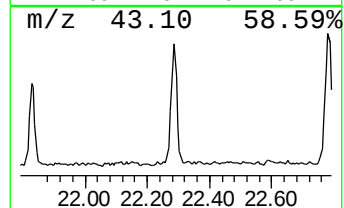
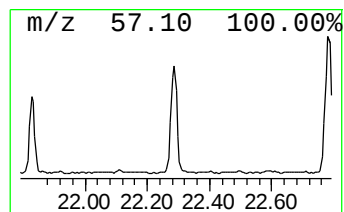
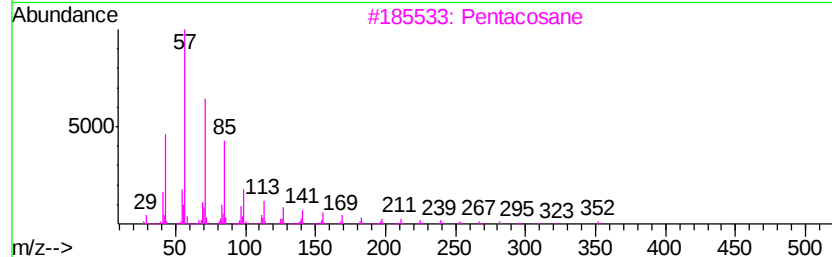
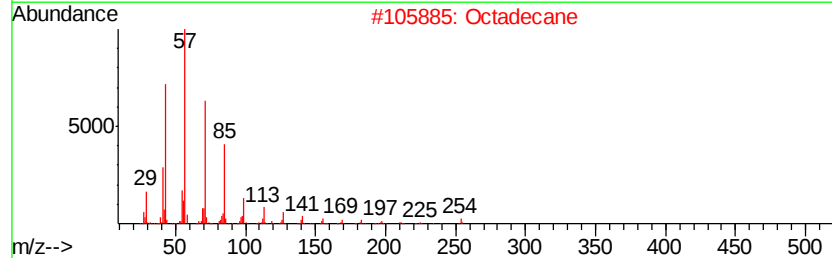
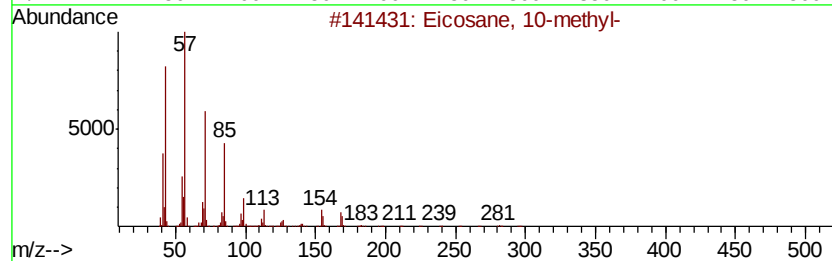
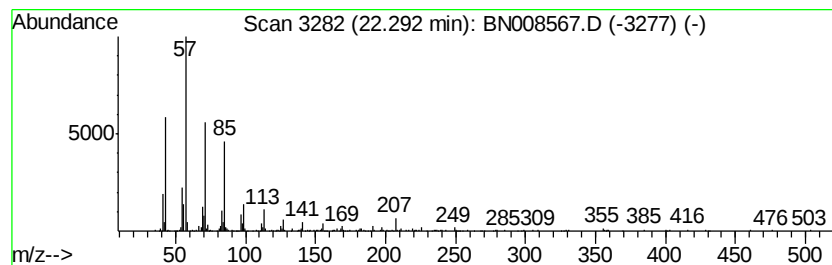
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.29	2.86 ng/ul	632134	Chrysene-d12	21.20

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane, 10-methyl-			296	C21H44	054833-23-7	87
2	Octadecane			254	C18H38	000593-45-3	83
3	Pentacosane			352	C25H52	000629-99-2	83
4	Tridecane, 2-methyl-			198	C14H30	001560-96-9	83
5	Docosane			310	C22H46	000629-97-0	83



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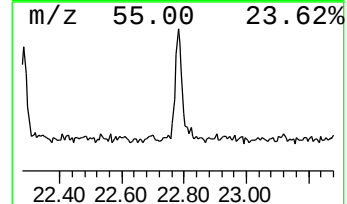
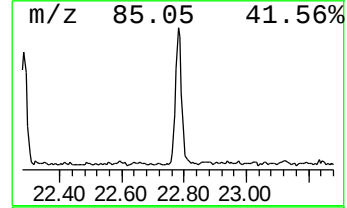
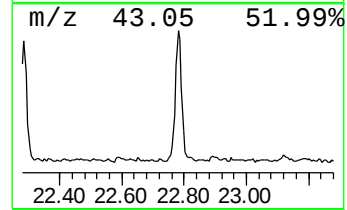
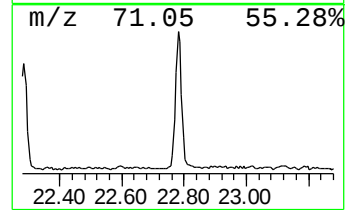
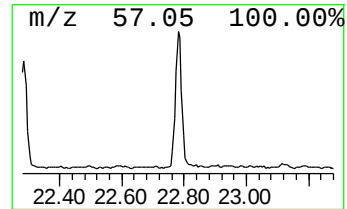
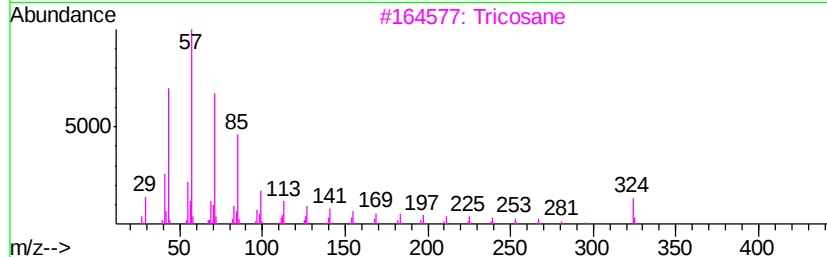
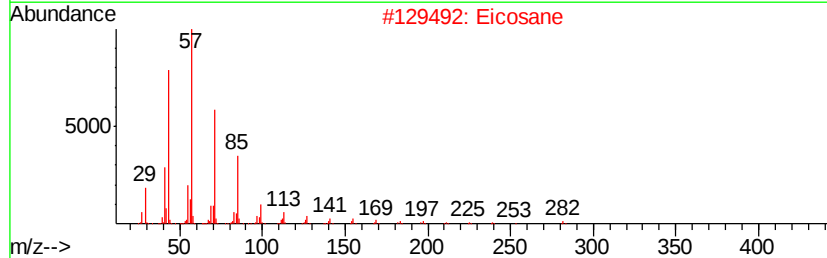
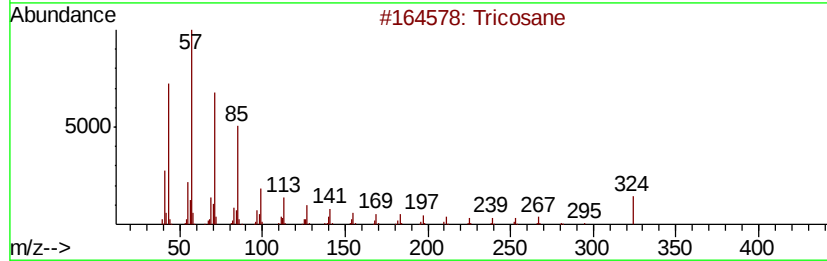
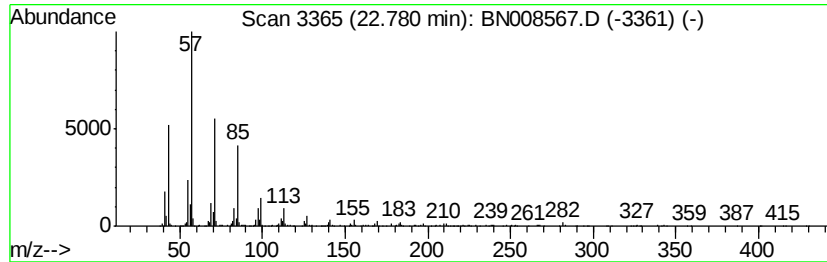
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.78	3.87 ng/ul	817476	Perylene-d12	23.37

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tricosane	324	C23H48	000638-67-5	94
2			Eicosane	282	C20H42	000112-95-8	93
3			Tricosane	324	C23H48	000638-67-5	91
4			Octacosane	394	C28H58	000630-02-4	90
5			2-methyloctacosane	408	C29H60	1000376-72-8	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111319\
Data File : BN008567.D
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Misc :
ALS Vial : 44 Sample Multiplier: 1

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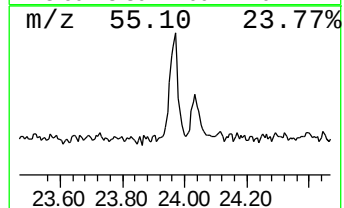
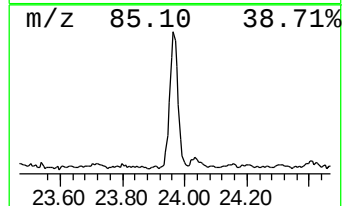
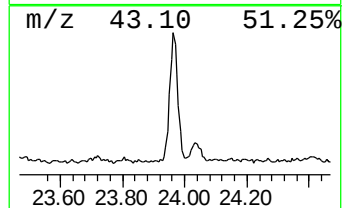
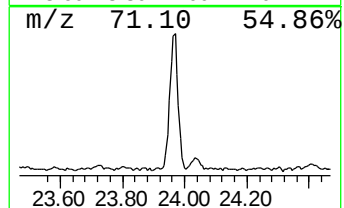
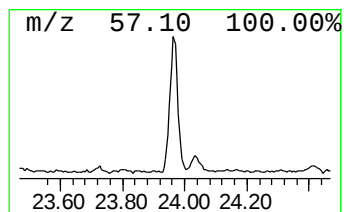
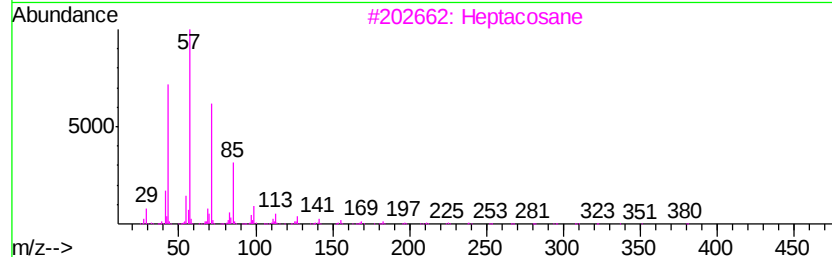
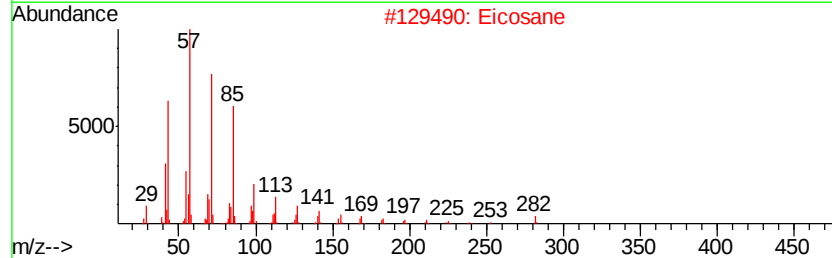
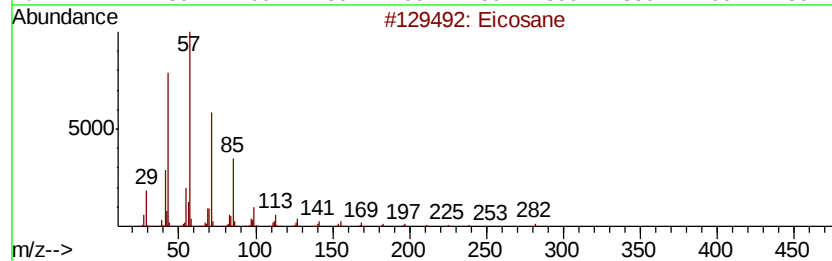
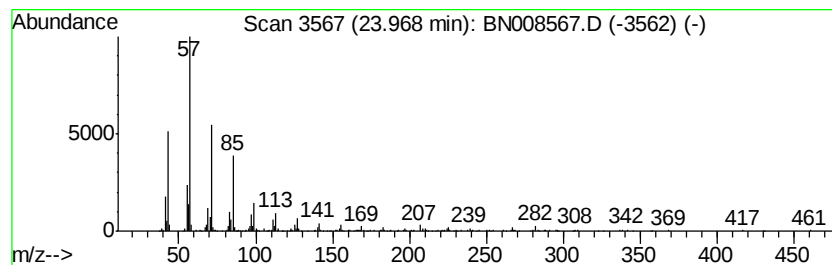
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN111319MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.97	4.30 ng/ul	906744	Perylene-d12	23.37

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane			282	C20H42	000112-95-8	93
2	Eicosane			282	C20H42	000112-95-8	91
3	Heptacosane			380	C27H56	000593-49-7	90
4	Nonadecane			268	C19H40	000629-92-5	90
5	Hentriacontane			437	C31H64	000630-04-6	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN111319\
Data File : BN008567.D
Acq On : 14 Nov 2019 21:31
Operator : JU
Sample : K5739-03
Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0CR1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN111319MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-chloro...	10.99	55.6	ng/ul	7355710	2	10.40	2644050	20.0
Benzene, 1-chloro...	11.20	13.8	ng/ul	1821280	2	10.40	2644050	20.0
4-Chloro-2-nitrop...	11.69	2.1	ng/ul	280154	2	10.40	2644050	20.0
Phenol, 5-chloro-...	12.37	2.2	ng/ul	399287	3	14.26	3649620	20.0
unknown-01	18.37	23.9	ng/ul	4928810	4	17.00	4118150	20.0
Octacosanol	20.95	4.1	ng/ul	910269	5	21.20	4426090	20.0
(DEL) Alkane: Str...	21.83	2.0	ng/ul	443715	5	21.20	4426090	20.0
(DEL) Alkane: Str...	22.29	2.9	ng/ul	632134	5	21.20	4426090	20.0
(DEL) Alkane: Str...	22.78	3.9	ng/ul	817476	6	23.37	4219350	20.0
(DEL) Alkane: Str...	23.97	4.3	ng/ul	906744	6	23.37	4219350	20.0