

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060319\
 Data File : BN006042.D
 Acq On : 03 Jun 2019 17:08
 Operator : JU/SJ
 Sample : K2924-21
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A42E8

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-BN060319MA.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	3.248	13	16	19	rBV	66461	90559	0.77%	0.033%
2	3.283	19	22	33	rVB	344056	470809	4.03%	0.174%
3	3.525	60	63	69	rBV2	18172	26691	0.23%	0.010%
4	3.695	88	92	98	rBV	59633	82582	0.71%	0.030%
5	3.972	135	139	145	rVB	21259	30973	0.26%	0.011%
6	4.283	187	192	211	rVB	4868082	6741591	57.68%	2.487%
7	4.872	286	292	303	rBV	1441106	2001091	17.12%	0.738%
8	5.477	390	395	401	rVB	43476	63424	0.54%	0.023%
9	5.930	468	472	478	rBV	36757	54287	0.46%	0.020%
10	6.877	626	633	636	rBV	2084445	3890020	33.28%	1.435%
11	6.930	639	642	656	rVV	2101684	3073654	26.30%	1.134%
12	7.066	659	665	680	rVB	1136542	1751669	14.99%	0.646%
13	7.266	692	699	701	rBV	1471855	2316202	19.82%	0.855%
14	7.295	701	704	712	rVB	1406399	2084231	17.83%	0.769%
15	7.730	771	778	785	rBV	1463709	2333533	19.96%	0.861%
16	7.789	785	788	795	rVB2	26362	44023	0.38%	0.016%
17	8.436	891	898	908	rBV	1720491	2815830	24.09%	1.039%
18	8.536	908	915	925	rVB	1349187	2136741	18.28%	0.788%
19	8.883	967	974	978	rBV	1324364	2248400	19.24%	0.830%
20	8.930	978	982	993	rVB	1113892	1776573	15.20%	0.656%
21	9.301	1040	1045	1051	rVB	51421	72178	0.62%	0.027%
22	9.607	1089	1097	1099	rBV	1135017	1843599	15.77%	0.680%
23	9.636	1099	1102	1112	rVB	1080901	1708749	14.62%	0.630%
24	10.142	1180	1188	1190	rBV	1845842	3096405	26.49%	1.142%
25	10.518	1243	1252	1256	rBV	2136675	3553763	30.40%	1.311%
26	10.571	1256	1261	1268	rVV	1635826	2784345	23.82%	1.027%
27	10.654	1268	1275	1288	rVV	1515110	2766889	23.67%	1.021%
28	10.854	1301	1309	1321	rVB	1680569	2716141	23.24%	1.002%
29	11.812	1464	1472	1489	rBV	3091531	5121605	43.82%	1.890%
30	11.948	1491	1495	1500	rVV4	15042	28163	0.24%	0.010%
31	12.054	1508	1513	1518	rVV	102527	164026	1.40%	0.061%
32	12.112	1518	1523	1530	rVV	28928	50673	0.43%	0.019%
33	12.412	1563	1574	1585	rBV	2693328	4384425	37.51%	1.618%
34	12.559	1592	1599	1608	rBV	1812609	2809105	24.03%	1.036%

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Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN060319MA.M
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35	12.806	1634	1641	1646	rVV	2996843	4641683	39.71%	1.713%
36	12.853	1646	1649	1666	rVV	262999	433773	3.71%	0.160%
37	12.989	1666	1672	1681	rVB6	20572	45133	0.39%	0.017%
38	13.212	1702	1710	1713	rBV	3621594	5531064	47.32%	2.041%
39	13.253	1713	1717	1729	rVB	1886434	3198073	27.36%	1.180%
40	13.795	1802	1809	1813	rVV	3121513	4439470	37.98%	1.638%
41	13.842	1813	1817	1825	rVV	3155865	4266022	36.50%	1.574%
42	13.965	1832	1838	1846	rVV	1812998	2543629	21.76%	0.939%
43	14.071	1848	1856	1873	rVB	3553884	5489121	46.96%	2.025%
44	14.383	1901	1909	1915	rBV2	2955930	4634286	39.65%	1.710%
45	14.448	1915	1920	1931	rVB	3386619	4798613	41.05%	1.771%
46	14.600	1935	1946	1966	rBV2	1961505	4186189	35.81%	1.545%
47	14.783	1971	1977	1994	rVB	2419560	3505209	29.99%	1.293%
48	15.012	2007	2016	2031	rBV	1874491	2834757	24.25%	1.046%
49	15.189	2041	2046	2050	rBV	49378	67400	0.58%	0.025%
50	15.242	2051	2055	2060	rVB2	33158	44600	0.38%	0.016%
51	15.383	2072	2079	2083	rBV	4498412	6421594	54.94%	2.369%
52	15.430	2083	2087	2095	rVV	2207296	2951375	25.25%	1.089%
53	15.524	2095	2103	2115	rVV3	2590721	4828728	41.31%	1.782%
54	15.618	2115	2119	2128	rVB3	74983	139783	1.20%	0.052%
55	16.106	2199	2202	2204	rBV3	25440	30518	0.26%	0.011%
56	16.171	2209	2213	2219	rVB5	22774	32799	0.28%	0.012%
57	16.442	2253	2259	2267	rBV	2031656	3031730	25.94%	1.119%
58	16.683	2295	2300	2304	rBV2	23887	34951	0.30%	0.013%
59	16.794	2313	2319	2334	rBV	2609047	3968120	33.95%	1.464%
60	17.141	2371	2378	2381	rVV	3680390	5372446	45.96%	1.982%
61	17.183	2381	2385	2389	rVV	3100984	4353219	37.24%	1.606%
62	17.236	2389	2394	2398	rVV2	4478548	7127608	60.98%	2.630%
63	17.271	2398	2400	2416	rVV	3410791	4415756	37.78%	1.629%
64	17.412	2419	2424	2433	rVV9	32759	98119	0.84%	0.036%
65	17.547	2440	2447	2453	rVV	2714821	4070309	34.82%	1.502%
66	17.618	2453	2459	2480	rVV	2815354	4068651	34.81%	1.501%
67	17.759	2480	2483	2486	rVV5	17652	27391	0.23%	0.010%
68	17.812	2489	2492	2496	rVB5	26311	30298	0.26%	0.011%
69	18.041	2527	2531	2536	rBV	311510	444513	3.80%	0.164%
70	18.112	2536	2543	2549	rVB	3103066	3932421	33.64%	1.451%
71	18.341	2578	2582	2585	rBV2	45455	63323	0.54%	0.023%

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72	18.377	2585	2588	2590	rBV2	21858	29592	0.25%	0.011%
73	18.565	2616	2620	2630	rVB5	50618	84590	0.72%	0.031%
74	18.918	2675	2680	2687	rBV2	108128	212303	1.82%	0.078%
75	18.977	2687	2690	2698	rVB3	75865	133585	1.14%	0.049%
76	19.212	2720	2730	2739	rBV	6758358	9376787	80.22%	3.460%
77	19.330	2747	2750	2757	rVB4	47594	67224	0.58%	0.025%
78	19.430	2764	2767	2769	rBV	41270	45272	0.39%	0.017%
79	19.547	2781	2787	2790	rBV2	5011651	7773869	66.51%	2.868%
80	19.577	2790	2792	2801	rVB	4924791	5609730	47.99%	2.070%
81	19.677	2806	2809	2815	rVB7	36849	66771	0.57%	0.025%
82	20.100	2878	2881	2885	rVB	83091	96480	0.83%	0.036%
83	20.412	2932	2934	2940	rVB7	25219	32695	0.28%	0.012%
84	20.488	2942	2947	2952	rBV	2967976	3306773	28.29%	1.220%
85	20.606	2962	2967	2970	rBV2	112646	143249	1.23%	0.053%
86	21.071	3042	3046	3058	rBV	773496	1302206	11.14%	0.480%
87	21.277	3076	3081	3086	rBV	4575450	4956583	42.40%	1.829%
88	21.347	3087	3093	3099	rVV2	5982194	11688829	100.00%	4.313%
89	21.400	3099	3102	3111	rVB	4139157	5052067	43.22%	1.864%
90	21.518	3118	3122	3132	rVB	954348	991279	8.48%	0.366%
91	21.965	3193	3198	3203	rBV	1723040	2098551	17.95%	0.774%
92	22.441	3273	3279	3285	rVB	2647437	3366632	28.80%	1.242%
93	22.959	3361	3367	3379	rBV2	5227475	8977242	76.80%	3.312%
94	23.512	3453	3461	3463	rBV2	2826547	5502197	47.07%	2.030%
95	23.547	3463	3467	3477	rVV2	4143807	8624503	73.78%	3.182%
96	23.653	3478	3485	3500	rVB	2138915	4157017	35.56%	1.534%
97	24.200	3571	3578	3588	rBV	2002796	3780958	32.35%	1.395%
98	24.959	3701	3707	3719	rVB	1072686	2394206	20.48%	0.883%
99	25.853	3852	3859	3867	rBV2	472078	1176758	10.07%	0.434%
100	25.965	3867	3878	3895	rVB2	2918897	8738001	74.76%	3.224%

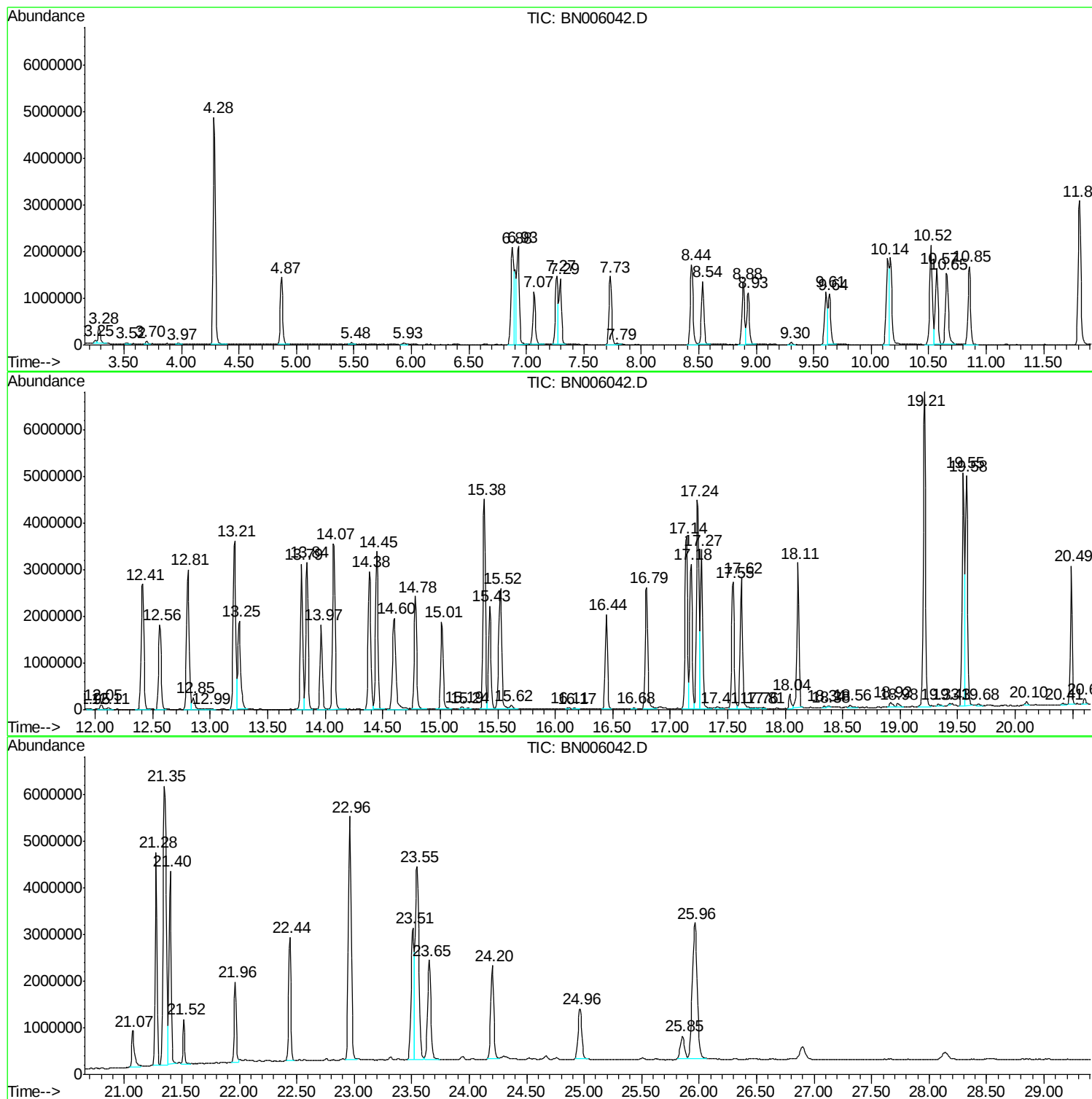
Sum of corrected areas: 271021572

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Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN060319MA.M
Quant Title : SVOA CALIBRATION

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



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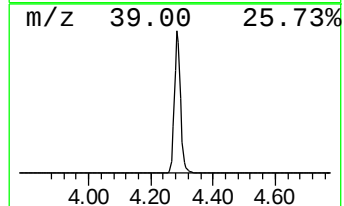
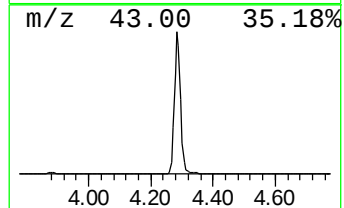
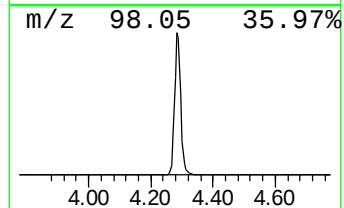
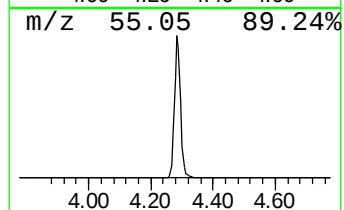
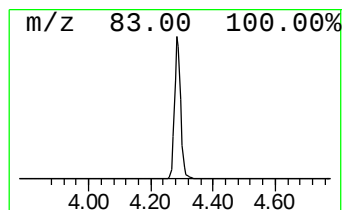
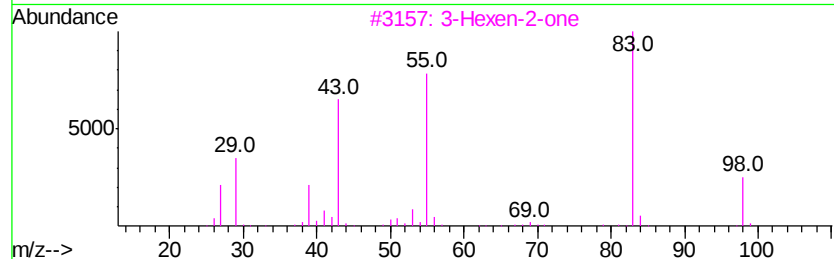
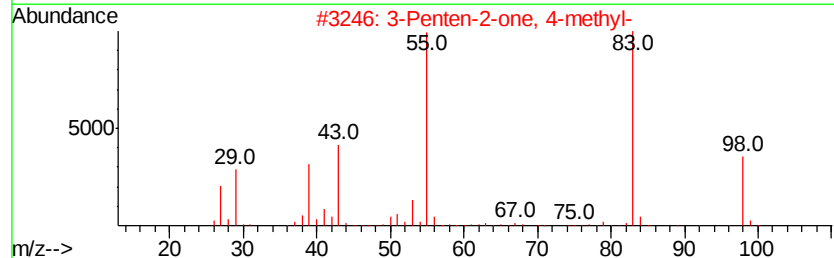
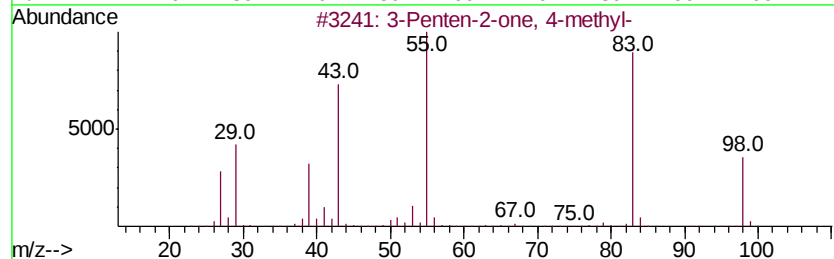
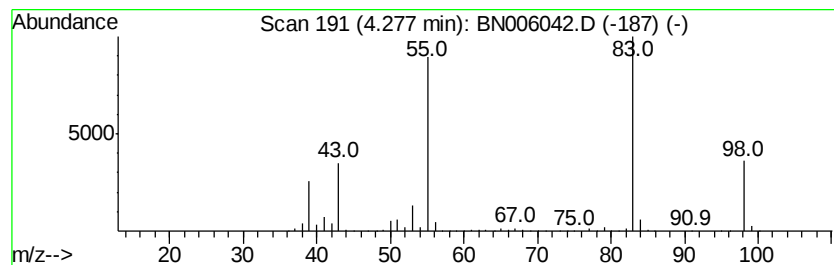
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN060319MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 3-Penten-2-one, 4-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.28	57.78 ng/ul	6741590	1,4-Dichlorobenzene-d4	7.73

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
2		3-Penten-2-one, 4-methyl-	98	C6H10O	000141-79-7	91
3		3-Hexen-2-one	98	C6H10O	000763-93-9	91
4		3-Hexen-2-one	98	C6H10O	000763-93-9	91
5		2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	90



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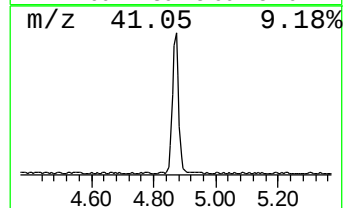
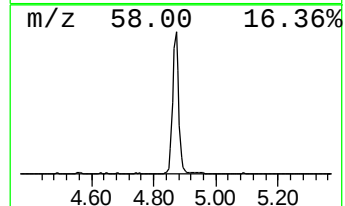
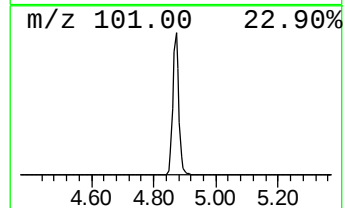
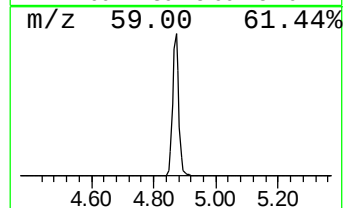
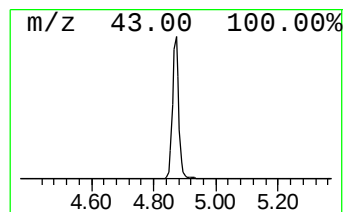
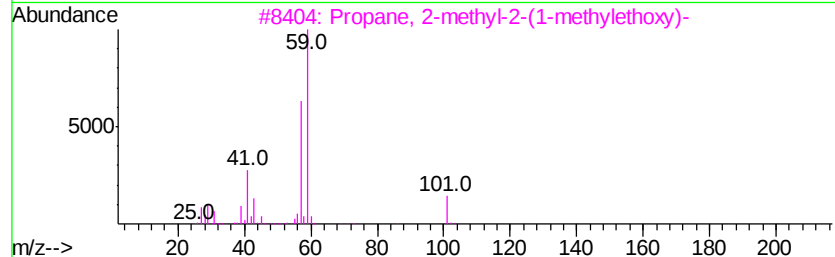
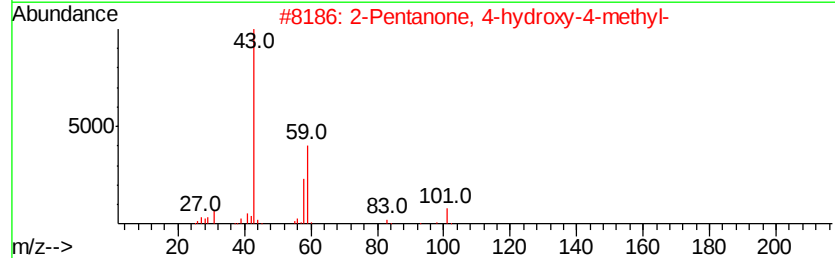
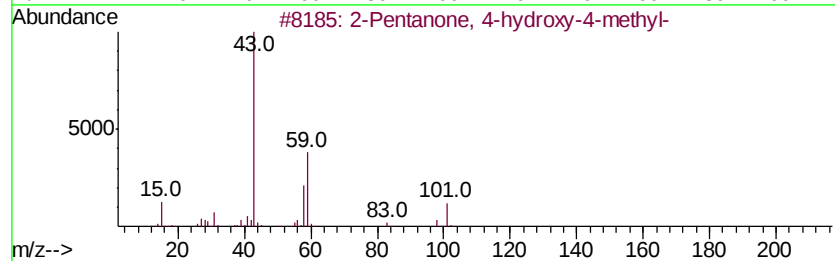
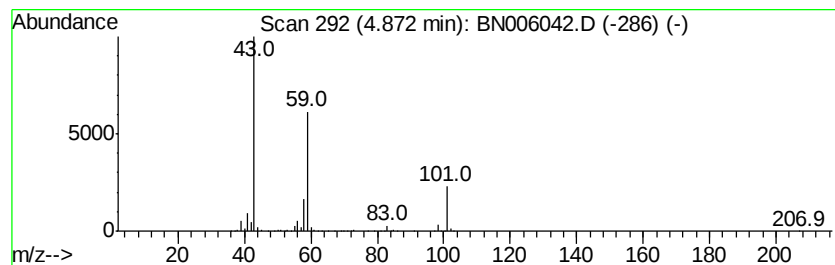
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.87	17.15 ng/ul	2001090	1,4-Dichlorobenzene-d4	7.73

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
3		Propane, 2-methyl-2-(1-methylethoxy)-	116	C7H16O	017348-59-3	53
4		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
5		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	23



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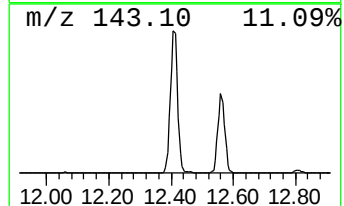
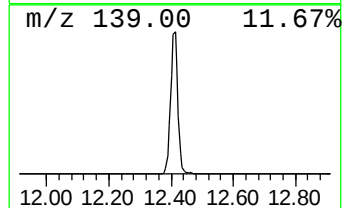
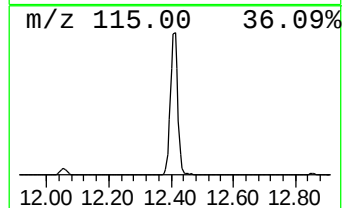
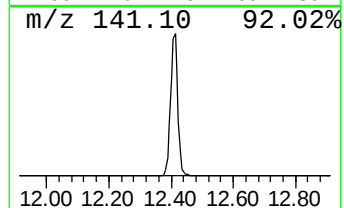
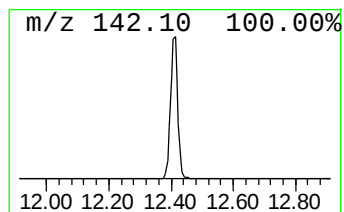
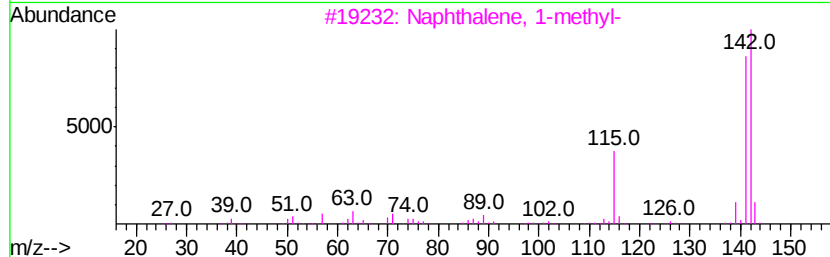
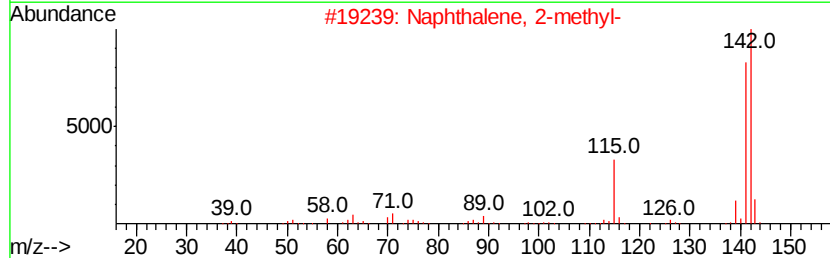
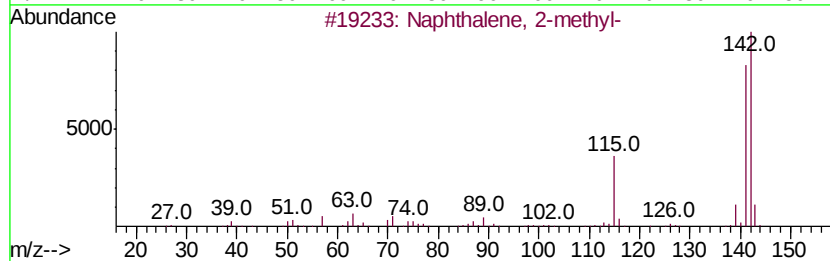
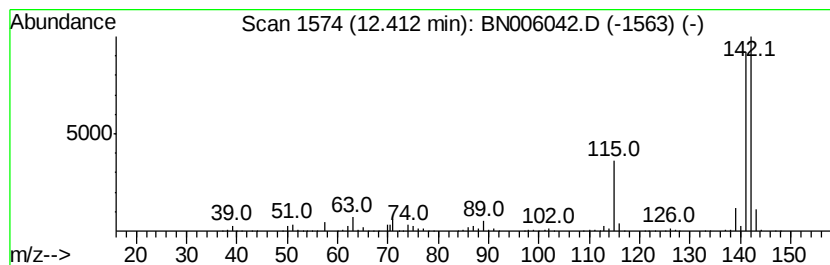
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Naphthalene, 1-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.41	24.67 ng/ul	4384430	Naphthalene-d8	10.52
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 2-methyl-	142 C11H10	000091-57-6 96
2		Naphthalene, 2-methyl-	142 C11H10	000091-57-6 96
3		Naphthalene, 1-methyl-	142 C11H10	000090-12-0 96
4		Naphthalene, 1-methyl-	142 C11H10	000090-12-0 94
5		Naphthalene, 1-methyl-	142 C11H10	000090-12-0 91



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Data File : BN006042.D
Acq On : 03 Jun 2019 17:08
Operator : JU/SJ
Sample : K2924-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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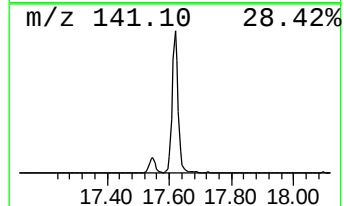
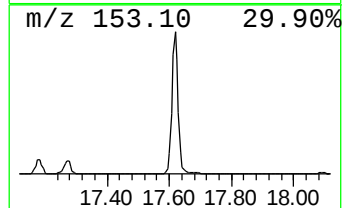
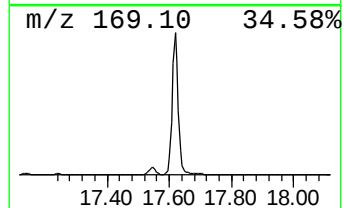
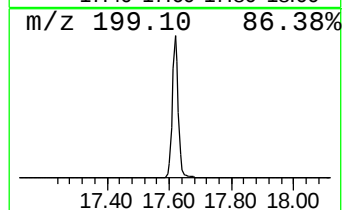
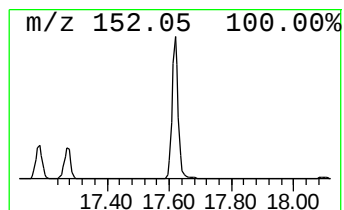
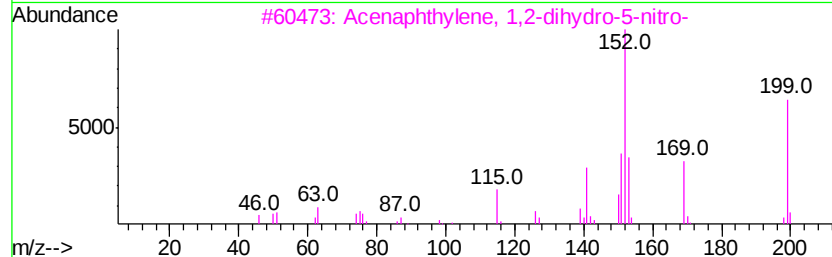
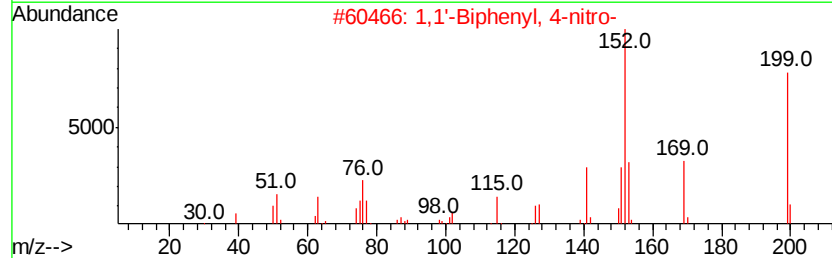
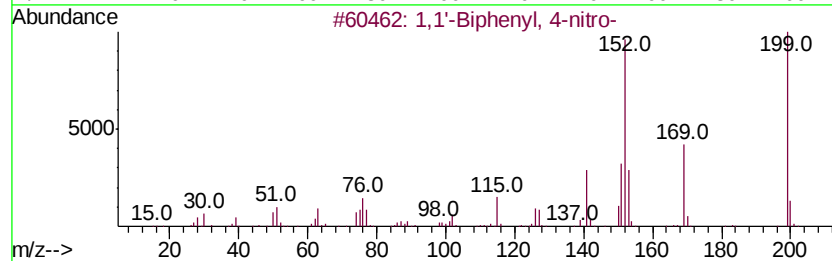
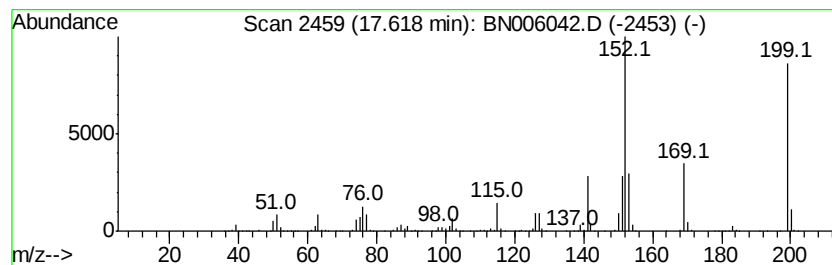
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1,1'-Biphenyl, 4-nitro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.62	15.15 ng/ul	4068650	Phenanthrene-d10	17.14

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1,1'-Biphenyl, 4-nitro-	199	C12H9NO2	000092-93-3	99
2		1,1'-Biphenyl, 4-nitro-	199	C12H9NO2	000092-93-3	98
3		Acenaphthylene, 1,2-dihydro-5-ni...	199	C12H9NO2	000602-87-9	94
4		1,1'-Biphenyl, 3-nitro-	199	C12H9NO2	002113-58-8	89
5		Acenaphthylene, 1,2-dihydro-5-ni...	199	C12H9NO2	000602-87-9	86



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Sample : K2924-21
Misc :
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Instrument :
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ClientSampleId :
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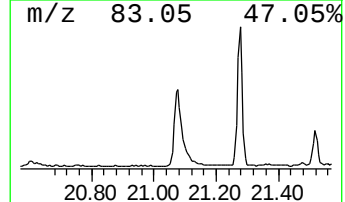
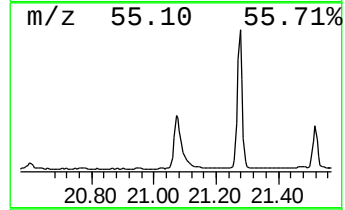
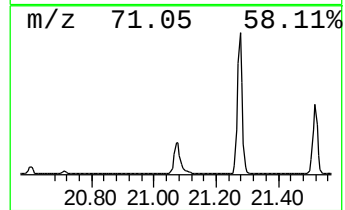
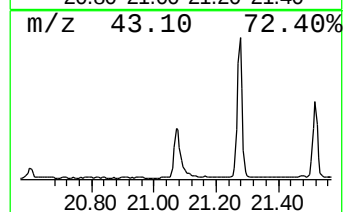
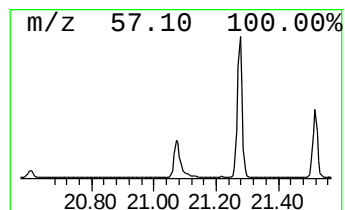
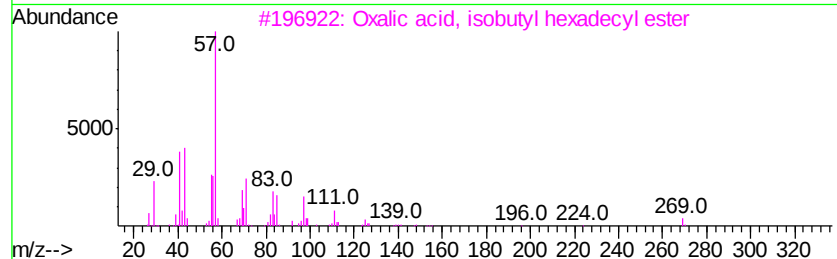
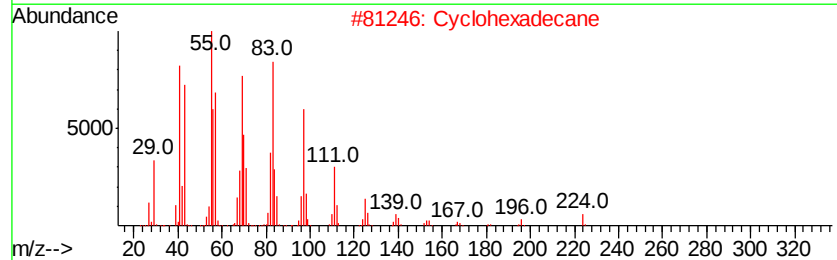
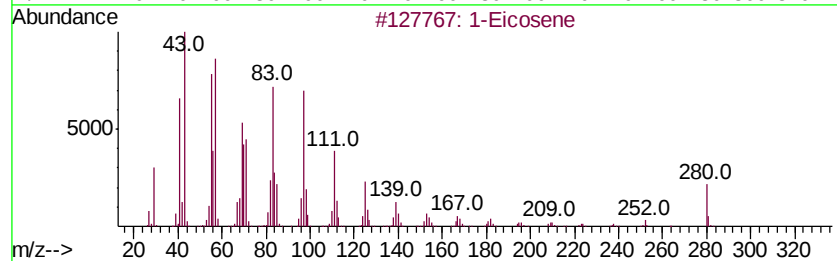
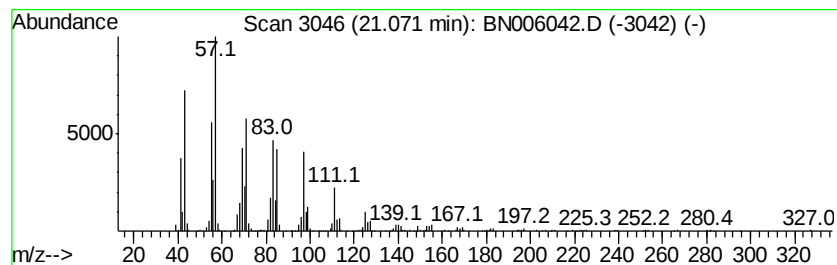
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Cyclic21.07 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.07	2.23 ng/ul	1302210	Chrysene-d12	21.36

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Eicosene		280	C20H40	003452-07-1	93
2	Cyclohexadecane		224	C16H32	000295-65-8	93
3	Oxalic acid, isobutyl hexadecyl ...		370	C22H42O4	1000309-38-1	91
4	Heptadecyl heptafluorobutyrate		452	C21H35F7O2	959085-66-6	76
5	Octadecyl trifluoroacetate		366	C20H37F3O2	079392-43-1	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060319\
Data File : BN006042.D
Acq On : 03 Jun 2019 17:08
Operator : JU/SJ
Sample : K2924-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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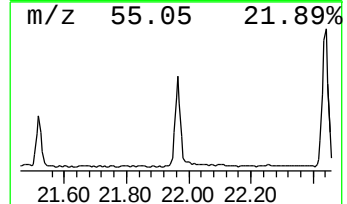
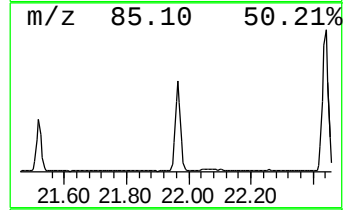
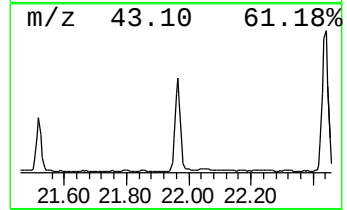
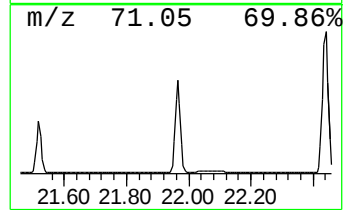
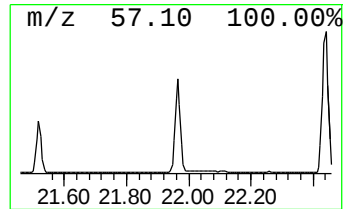
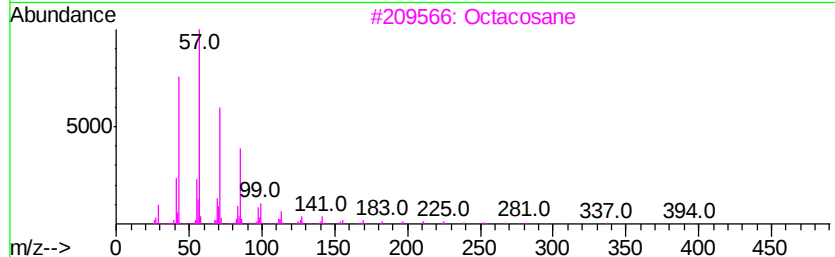
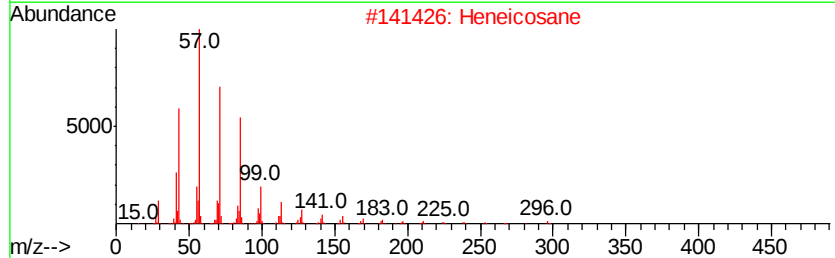
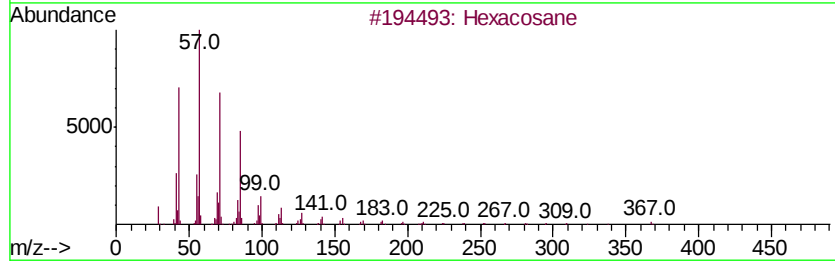
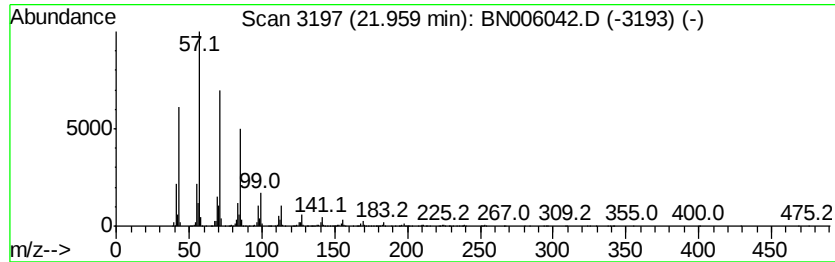
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.96	3.59 ng/ul	2098550	Chrysene-d12	21.36

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	94
2			Heneicosane	296	C21H44	000629-94-7	91
3			Octacosane	394	C28H58	000630-02-4	91
4			Heneicosane, 11-(1-ethylpropyl)-	366	C26H54	055282-11-6	91
5			Heneicosane	296	C21H44	000629-94-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060319\
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Acq On : 03 Jun 2019 17:08
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Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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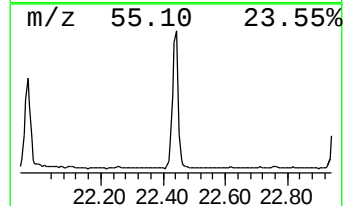
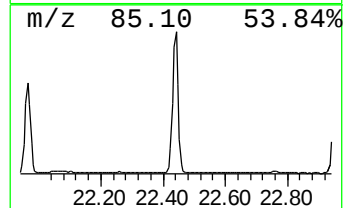
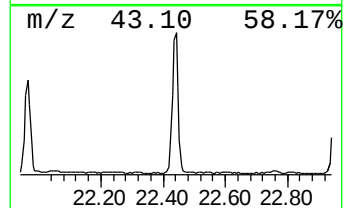
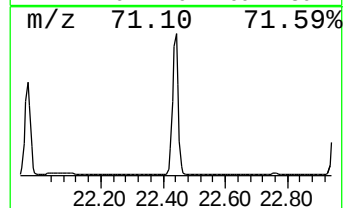
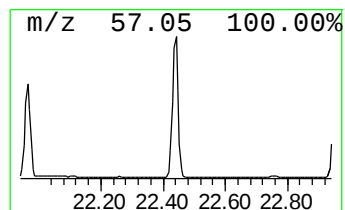
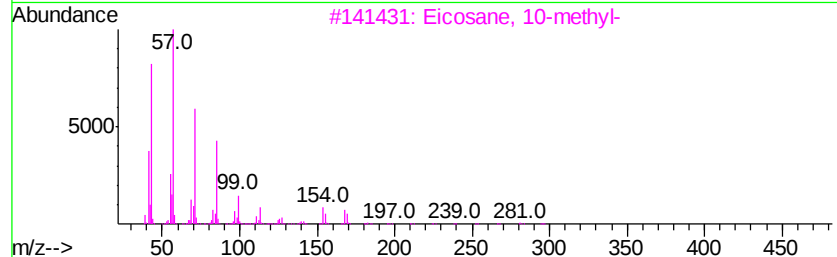
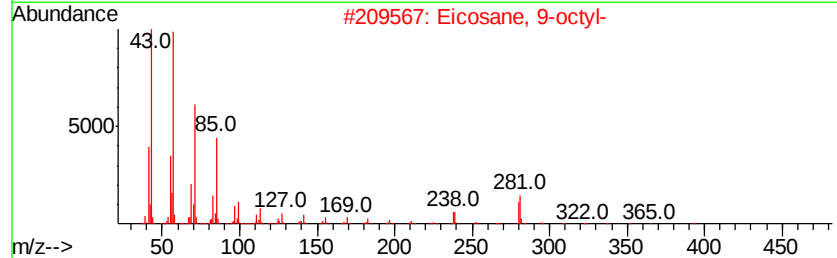
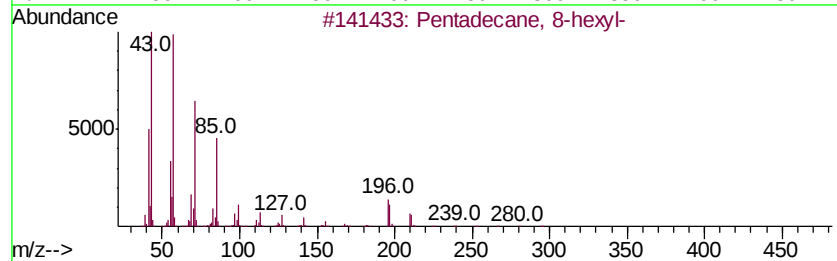
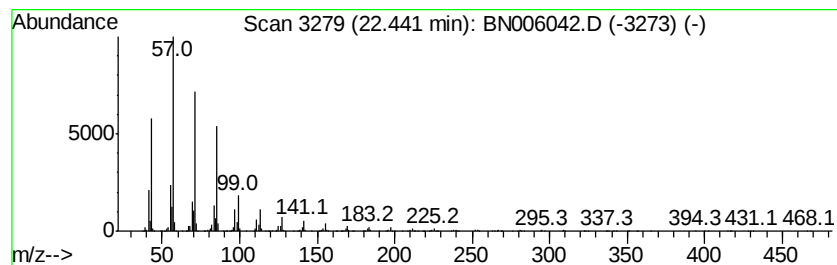
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.44	5.76 ng/ul	3366630	Chrysene-d12	21.36

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	94
2		Eicosane, 9-octyl-	394	C28H58	013475-77-9	94
3		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
4		Heneicosane	296	C21H44	000629-94-7	91
5		2-methyloctacosane	408	C29H60	1000376-72-8	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060319\
Data File : BN006042.D
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Sample : K2924-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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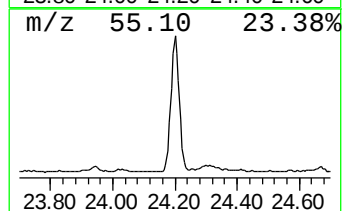
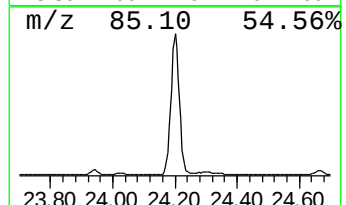
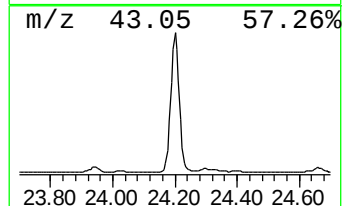
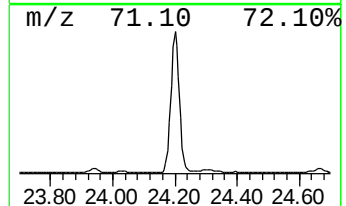
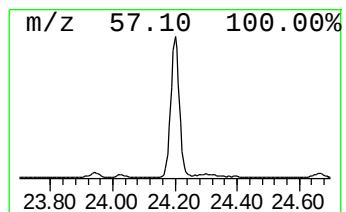
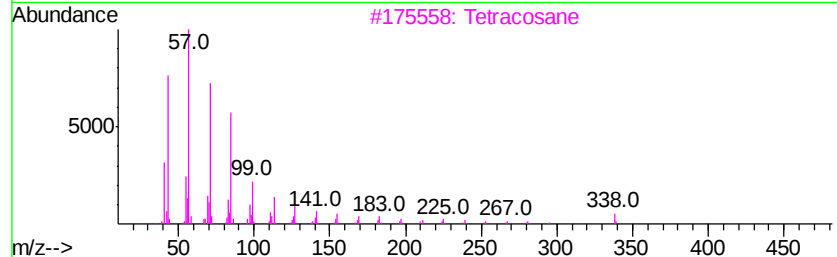
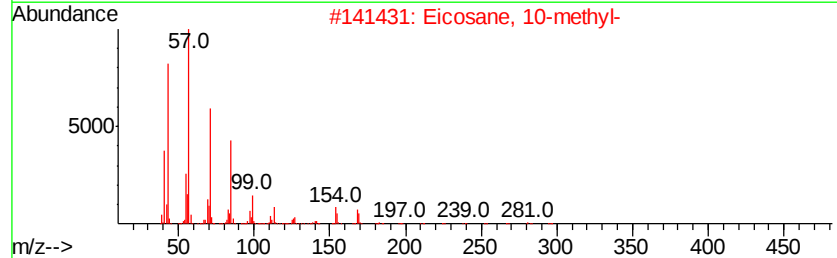
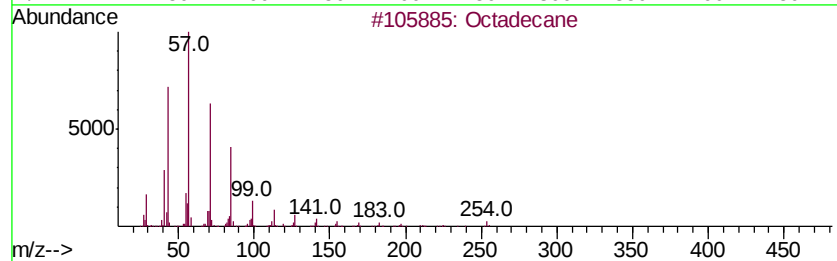
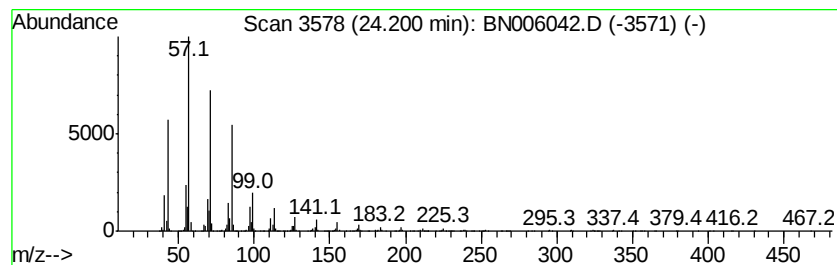
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.20	18.19 ng/ul	3780960	Perylene-d12	23.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	93
2		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
3		Tetracosane	338	C24H50	000646-31-1	91
4		Hexacosane	366	C26H54	000630-01-3	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060319\
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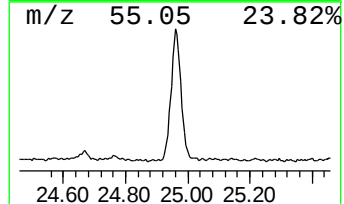
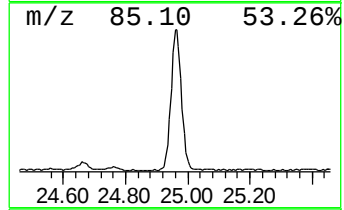
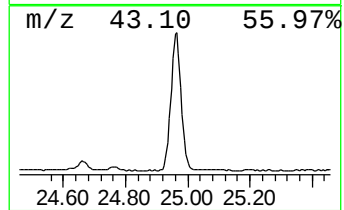
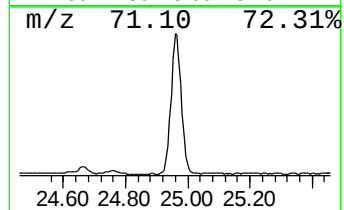
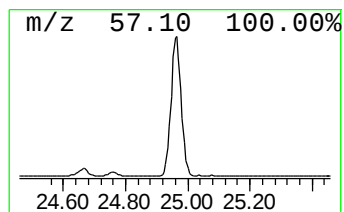
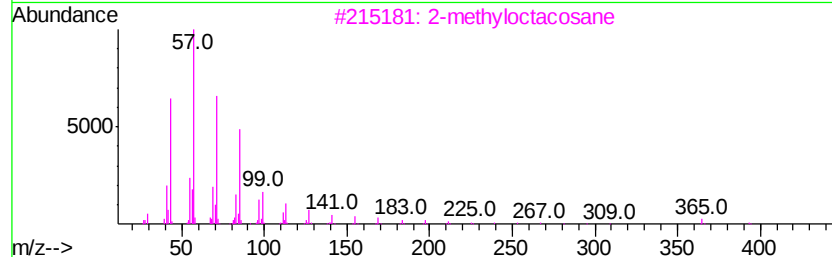
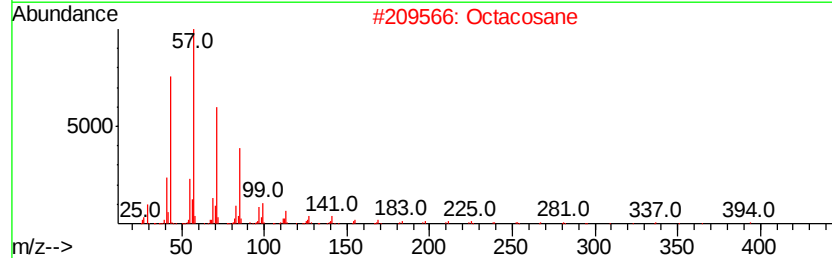
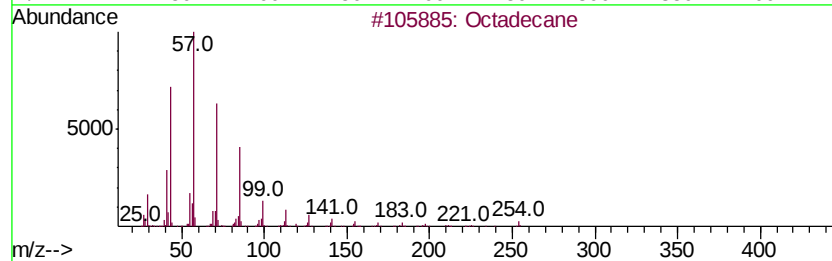
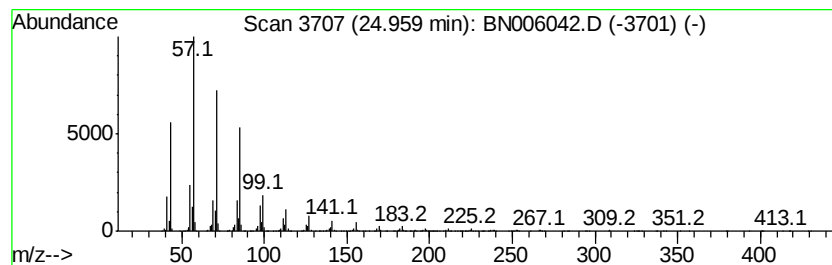
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.96	11.52 ng/ul	2394210	Perylene-d12	23.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Octadecane	254	C18H38	000593-45-3	96
2		Octacosane	394	C28H58	000630-02-4	91
3		2-methyloctacosane	408	C29H60	1000376-72-8	91
4		Eicosane, 10-methyl-	296	C21H44	054833-23-7	91
5		Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060319\
Data File : BN006042.D
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Misc :
ALS Vial : 9 Sample Multiplier: 1

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ClientSampleId :
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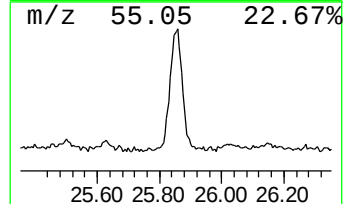
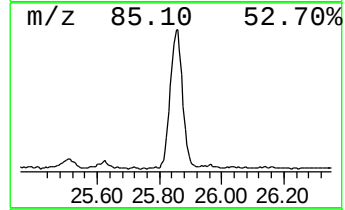
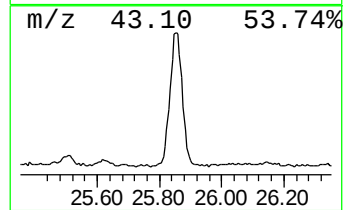
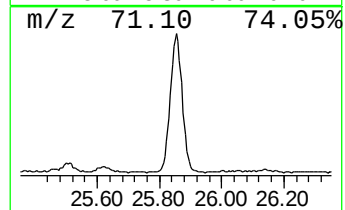
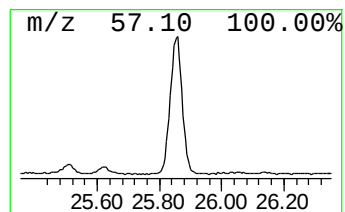
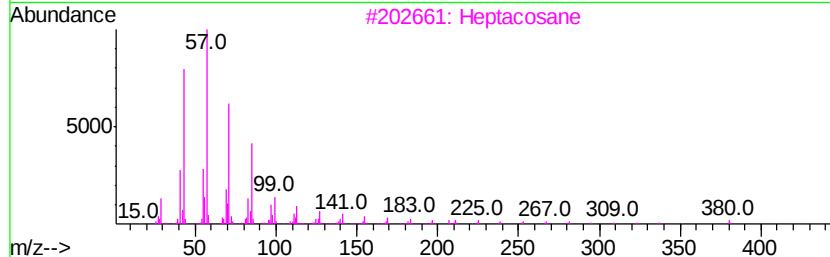
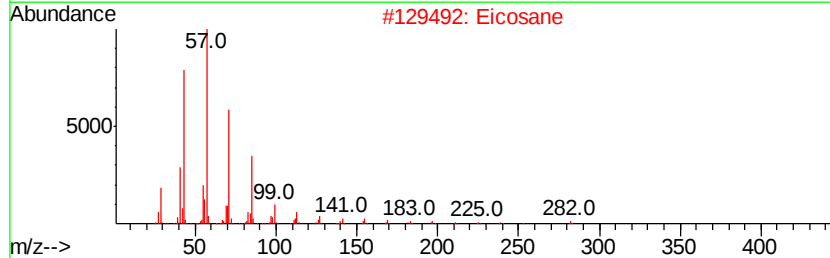
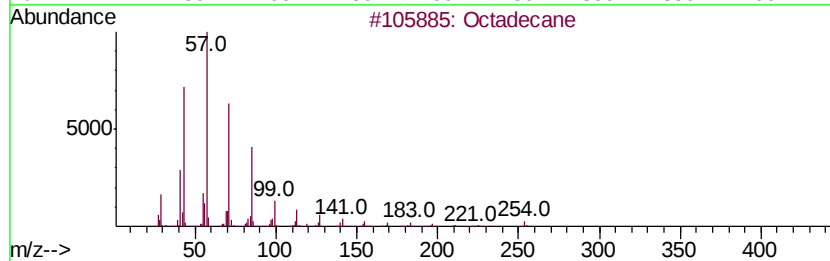
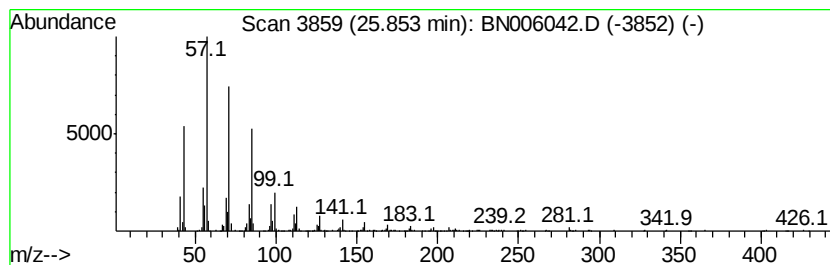
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN060319MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.85	5.66 ng/ul	1176760	Perylene-d12	23.65

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	95
2			Eicosane	282	C20H42	000112-95-8	95
3			Heptacosane	380	C27H56	000593-49-7	91
4			Heneicosane	296	C21H44	000629-94-7	91
5			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN060319\
Data File : BN006042.D
Acq On : 03 Jun 2019 17:08
Operator : JU/SJ
Sample : K2924-21
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
A42E8

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN060319MA.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
3-Penten-2-one, 4...	4.28	57.8	ng/ul	6741590	1	7.73	2333530	20.0
2-Pentanone, 4-hy...	4.87	17.1	ng/ul	2001090	1	7.73	2333530	20.0
Naphthalene, 1-me...	12.41	24.7	ng/ul	4384430	2	10.52	3553760	20.0
1,1'-Biphenyl, 4-...	17.62	15.2	ng/ul	4068650	4	17.14	5372450	20.0
(DEL) Alkane: Cyc...	21.07	2.2	ng/ul	1302210	5	21.36	11688800	20.0
(DEL) Alkane: Str...	21.96	3.6	ng/ul	2098550	5	21.36	11688800	20.0
(DEL) Alkane: Str...	22.44	5.8	ng/ul	3366630	5	21.36	11688800	20.0
(DEL) Alkane: Str...	24.20	18.2	ng/ul	3780960	6	23.65	4157020	20.0
(DEL) Alkane: Str...	24.96	11.5	ng/ul	2394210	6	23.65	4157020	20.0
(DEL) Alkane: Str...	25.85	5.7	ng/ul	1176760	6	23.65	4157020	20.0