LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102419\
Data File : BM023397.D
Acq On : 25 Oct 2019 06:00
Operator : JU
Sample : K5403-07
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFB96

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.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	3.046	8	10	13	rVB2	66987	61167	0.34%	0.065%
2	3.087	13	17	21	rVB	1372339	1342721	7.51%	1.428%
3	3.393	65	69	70	rBV2	91861	100355	0.56%	0.107%
4	3.416	70	73	79	rVB3	124343	153730	0.86%	0.163%
5	3.875	146	151	157	rVB	606255	769125	4.30%	0.818%
6	3.993	167	171	175	rVB2	61798	73661	0.41%	0.078%
7	4.516	257	260	265	rVB	47729	61755	0.35%	0.066%
8	4.622	275	278	282	rVB2	46087	60350	0.34%	0.064%
9	4.710	290	293	298	rVB	79527	103737	0.58%	0.110%
10	4.975	335	338	344	rVB	44103	61205	0.34%	0.065%
11	5.169	368	371	378	rVB2	42569	66593	0.37%	0.071%
12	5.669	451	456	465	rVB6	33192	81711	0.46%	0.087%
13	5.934	494	501	508	rBV7	40433	92180	0.52%	0.098%
14	6.140	530	536	545	rVB	1831936	2676296	14.98%	2.846%
15	6.216	545	549	555	rBV2	43232	63845	0.36%	0.068%
16	6.298	559	563	566	rVV5	29057	59071	0.33%	0.063%
17	6.545	600	605	611	rVV5	23303	57095	0.32%	0.061%
18	6.651	620	623	628	rVB	60947	90052	0.50%	0.096%
19	6.793	642	647	656	rVV7	51859	118655	0.66%	0.126%
20	6.987	674	680	684	rBV5	20527	43673	0.24%	0.046%
21	7.287	727	731	737	rVB5	39250	70195	0.39%	0.075%
22	7.463	756	761	764	rBV4	21667	42083	0.24%	0.045%
23	7.545	769	775	778	rBV2	41451	77795	0.44%	0.083%
24	7.634	783	790	800	rVB	2295037	3648529	20.42%	3.880%
25	7.881	828	832	834	rBV4	33379	41229	0.23%	0.044%
26	7.998	848	852	859	rVB5	31885	60778	0.34%	0.065%
27	8.192	880	885	890	rVB3	60568	110406	0.62%	0.117%
28	8.322	903	907	914	rVB	39757	71240	0.40%	0.076%
29	8.457	926	930	934	rVB7	29512	51790	0.29%	0.055%
30	8.522	934	941	945	rBV5	31929	64427	0.36%	0.069%
31	8.745	971	979	988	rVB7	60673	152137	0.85%	0.162%
32	8.845	988	996	997	rBV7	28969	49970	0.28%	0.053%
33	8.881	999	1002	1011	rVB3	55870	104495	0.58%	0.111%
34	8.969	1011	1017	1018	rBV5	33433	56479	0.32%	0.060%

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Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
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Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Title : SVOA CALIBRATION

35	8.998	1018	1022	1028	rVB7	50769	97442	0.55%	0.104%
36	9.075	1028	1035	1037	rBV7	20847	44116	0.25%	0.047%
37	9.228	1056	1061	1064	rBV7	22844	43415	0.24%	0.046%
38	9.292	1069	1072	1075	rVV3	41397	61131	0.34%	0.065%
39	9.345	1077	1081	1089	rVV7	45503	88994	0.50%	0.095%
40	9.510	1103	1109	1115	rBV6	24206	60120	0.34%	0.064%
41	9.563	1115	1118	1122	rVV5	30052	51519	0.29%	0.055%
42	9.610	1122	1126	1134	rVB9	41569	97983	0.55%	0.104%
43	10.028	1192	1197	1207	rVB9	22691	54688	0.31%	0.058%
44	10.133	1207	1215	1218	rBV9	26971	54136	0.30%	0.058%
45	10.292	1237	1242	1246	rBV7	25513	44213	0.25%	0.047%
46	10.416	1253	1263	1269	rBV	2988219	5096299	28.52%	5.419%
47	10.463	1269	1271	1276	rVB	91547	128215	0.72%	0.136%
48	10.557	1282	1287	1291	rVB6	26311	40341	0.23%	0.043%
49	10.610	1291	1296	1300	rBV3	54785	84038	0.47%	0.089%
50	10.833	1329	1334	1340	rBV9	23924	45701	0.26%	0.049%
51	10.922	1340	1349	1353	rVV10	29046	68418	0.38%	0.073%
52	11.110	1378	1381	1385	rVV5	32035	57927	0.32%	0.062%
53	11.280	1404	1410	1416	rBV9	21609	46130	0.26%	0.049%
54	11.475	1437	1443	1447	rBV2	90612	178786	1.00%	0.190%
55	11.880	1507	1512	1520	rVB5	42514	89756	0.50%	0.095%
56	12.086	1543	1547	1558	rVB2	99029	193630	1.08%	0.206%
57	12.169	1558	1561	1565	rBV6	28564	54694	0.31%	0.058%
58	12.310	1580	1585	1591	rBV	112541	169947	0.95%	0.181%
59	12.722	1648	1655	1660	rVB8	45093	99577	0.56%	0.106%
60	12.792	1662	1667	1671	rVV7	29491	61931	0.35%	0.066%
61	12.839	1671	1675	1680	rVV	126073	179392	1.00%	0.191%
62	12.886	1680	1683	1689	rVB6	29162	48948	0.27%	0.052%
63	13.069	1709	1714	1717	rBV4	60090	105870	0.59%	0.113%
64	13.104	1718	1720	1726	rVB4	34653	70221	0.39%	0.075%
65	13.222	1737	1740	1744	rVB5	34751	46303	0.26%	0.049%
66	13.310	1751	1755	1758	rBV4	31912	60298	0.34%	0.064%
67	13.445	1774	1778	1785	rBV7	45543	106721	0.60%	0.113%
68	13.610	1801	1806	1810	rVV2	151160	223534	1.25%	0.238%
69	13.663	1810	1815	1819	rVV3	101527	164476	0.92%	0.175%
70	13.716	1819	1824	1828	rVB2	151920	217688	1.22%	0.231%
71	13.792	1832	1837	1841	rVV2	273803	437501	2.45%	0.465%

LSC Area Percent Report

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Misc :
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Instrument : BNA_M
ClientSampleId : BFB96

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
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Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Title : SVOA CALIBRATION

72	13.839	1842	1845	1850	rVB3	70345	106463	0.60%	0.113%
73	14.010	1870	1874	1876	rBV4	60611	87540	0.49%	0.093%
74	14.121	1890	1893	1898	rVB2	124433	165052	0.92%	0.176%
75	14.210	1903	1908	1912	rBV6	48750	94195	0.53%	0.100%
76	14.286	1912	1921	1928	rBV	4020087	6371146	35.65%	6.775%
77	14.492	1954	1956	1961	rBV4	31702	62263	0.35%	0.066%
78	14.651	1980	1983	1986	rBV4	47288	77039	0.43%	0.082%
79	14.692	1986	1990	1998	rVB5	91205	210697	1.18%	0.224%
80	14.798	2005	2008	2011	rBV	120570	150307	0.84%	0.160%
81	14.833	2012	2014	2021	rVB5	109891	164726	0.92%	0.175%
82	15.086	2053	2057	2065	rVB4	130796	207567	1.16%	0.221%
83	15.168	2068	2071	2079	rBV7	53281	102651	0.57%	0.109%
84	15.333	2096	2099	2103	rBV	140347	200466	1.12%	0.213%
85	15.580	2137	2141	2146	rVB	193232	290204	1.62%	0.309%
86	16.027	2214	2217	2218	rBV3	115202	116773	0.65%	0.124%
87	16.057	2219	2222	2229	rVB	438215	597407	3.34%	0.635%
88	16.186	2240	2244	2248	rVB	327962	424548	2.38%	0.451%
89	16.345	2266	2271	2277	rBV	8358777	11084375	62.03%	11.786%
90	16.704	2328	2332	2339	rVB	856803	1067028	5.97%	1.135%
91	16.880	2358	2362	2366	rVB2	331232	448449	2.51%	0.477%
92	17.033	2382	2388	2392	rBV2	3692287	5615137	31.42%	5.971%
93	17.074	2392	2395	2400	rVB	958552	1279623	7.16%	1.361%
94	19.098	2735	2739	2746	rVB	1380895	1863753	10.43%	1.982%
95	19.162	2747	2750	2758	rVB	492400	701510	3.93%	0.746%
96	21.227	3097	3101	3105	rVB	3419979	4213595	23.58%	4.480%
97	21.503	3143	3148	3151	rBV	14818210	17870679	100.00%	19.002%
98	21.721	3181	3185	3192	rBV	6059220	7970717	44.60%	8.475%
99	23.427	3471	3475	3487	rVB	3275618	6029870	33.74%	6.412%
100	27.132	4099	4105	4119	rVB2	2215684	6959030	38.94%	7.400%

Sum of corrected areas: 94045439

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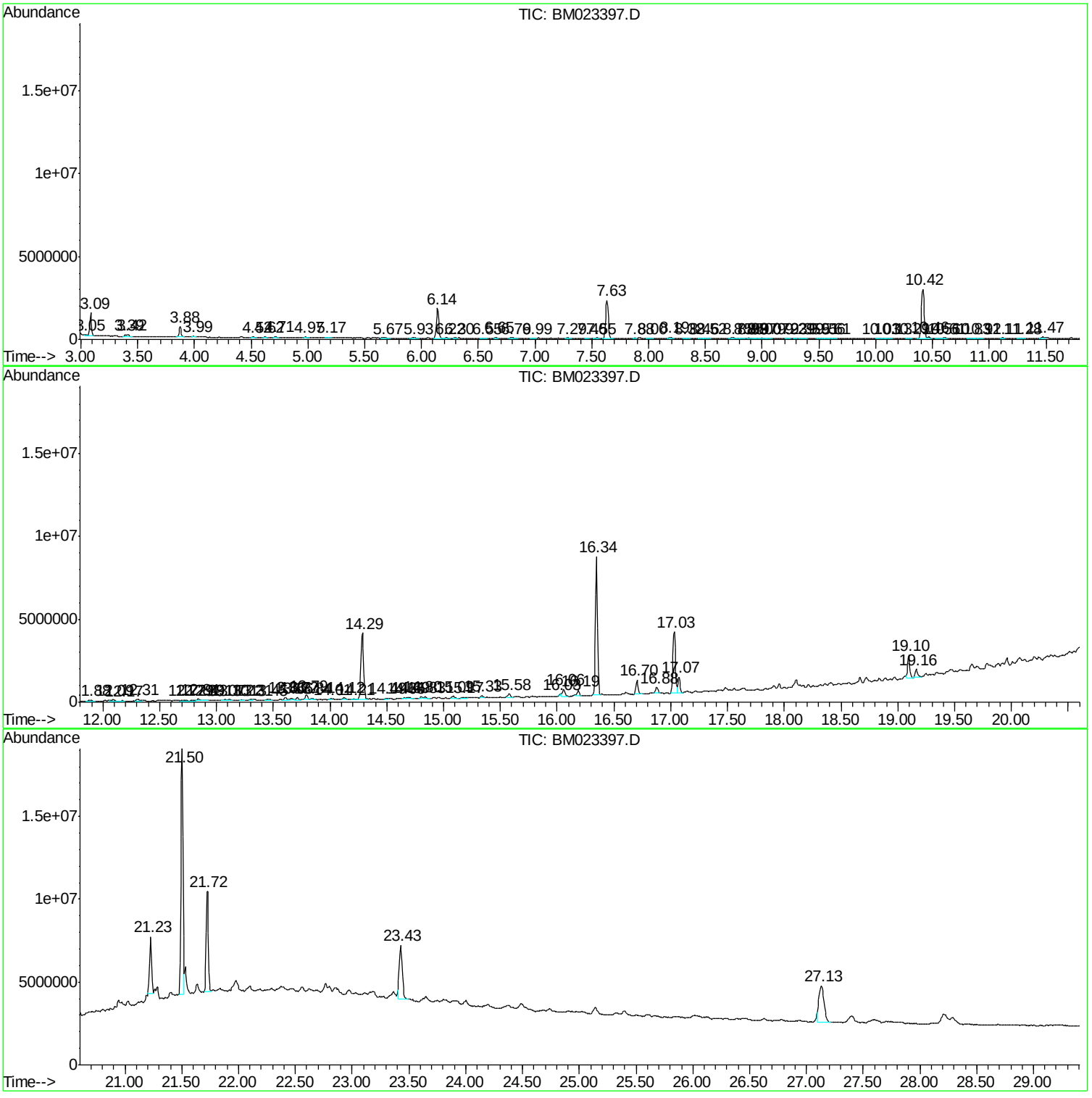
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
BFB96

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

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



Library Search Compound Report

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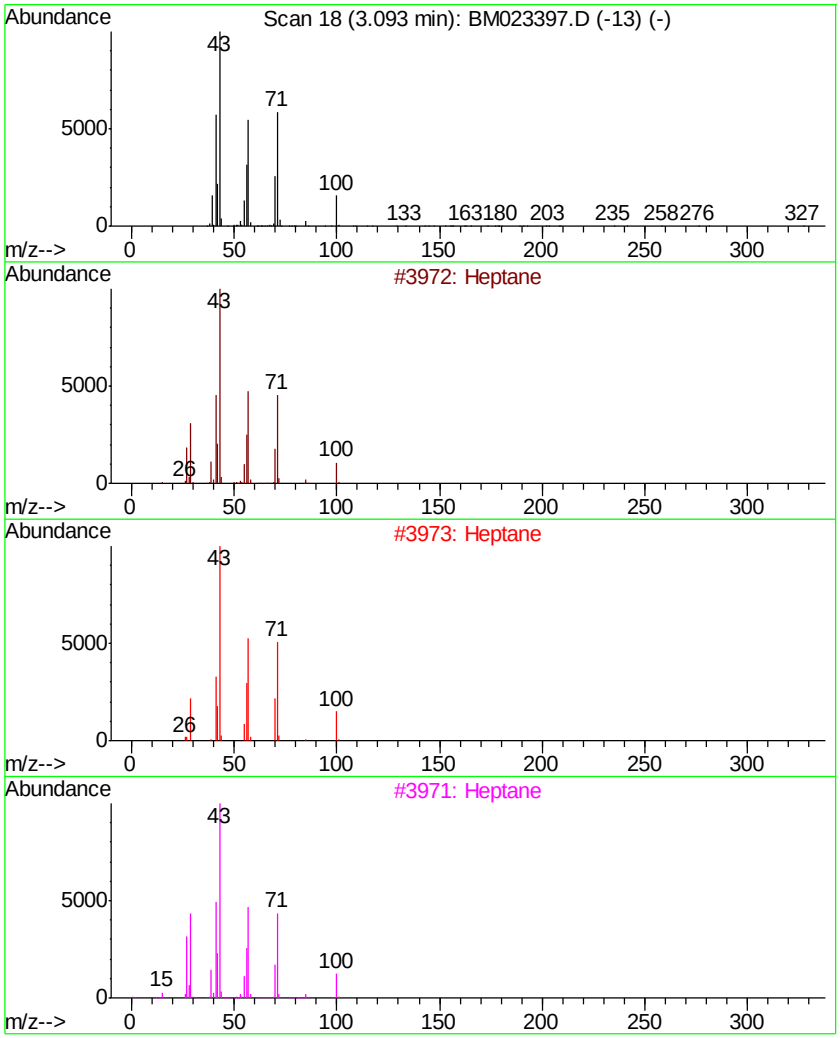
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

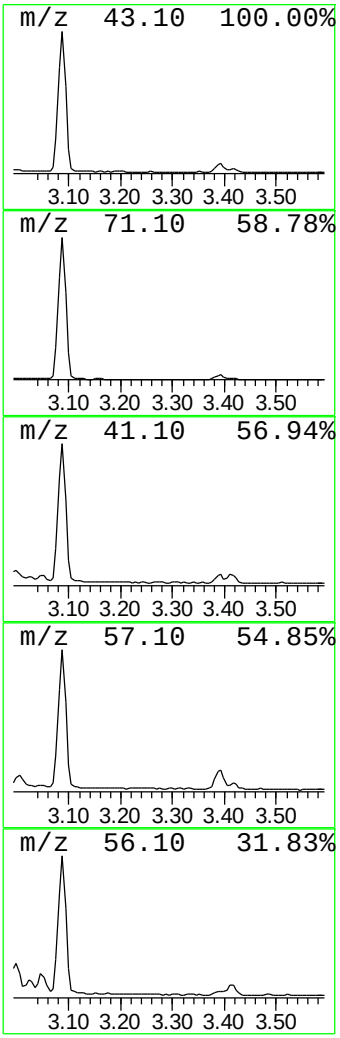
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.09	7.36 ng/ul	1342720	1,4-Dichlorobenzene-d4	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane	100	C7H16	000142-82-5	95
2	Heptane	100	C7H16	000142-82-5	91
3	Heptane	100	C7H16	000142-82-5	90
4	Heptane	100	C7H16	000142-82-5	87
5	Oxalic acid, isobutyl pentyl ester	216	C11H20O4	1000309-37-0	59





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Data File : BM023397.D
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Instrument : BNA_M
ClientSampleId : BFB96

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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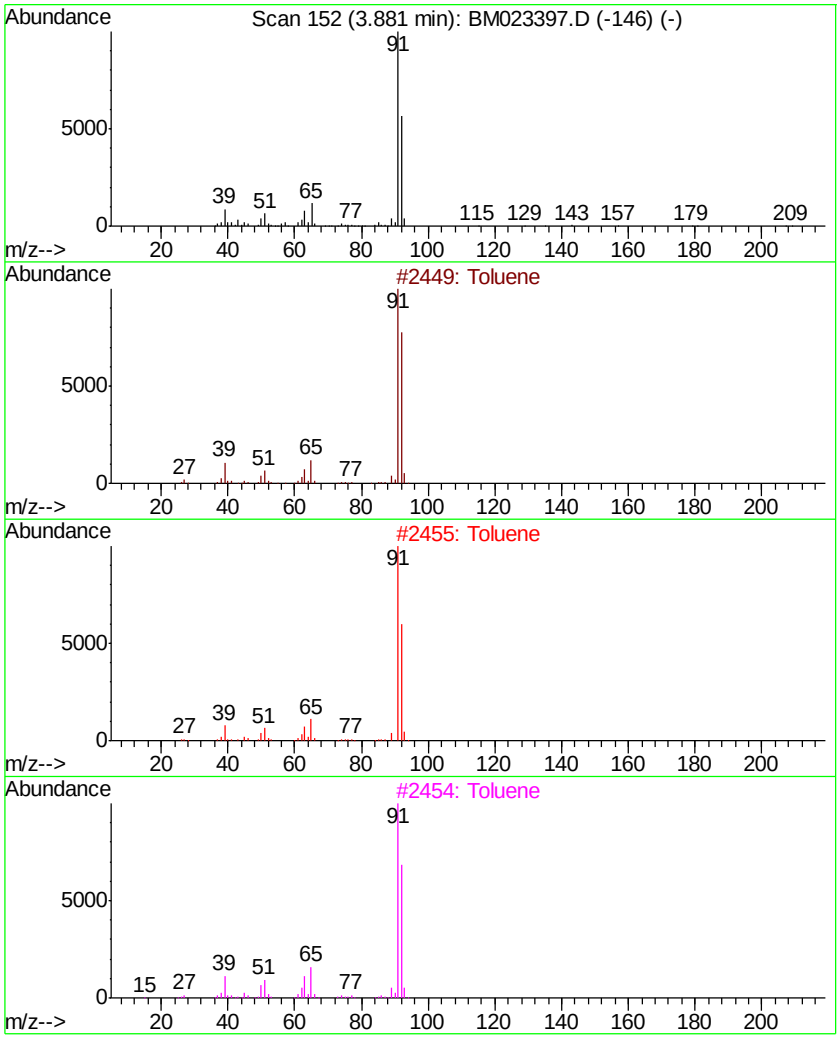
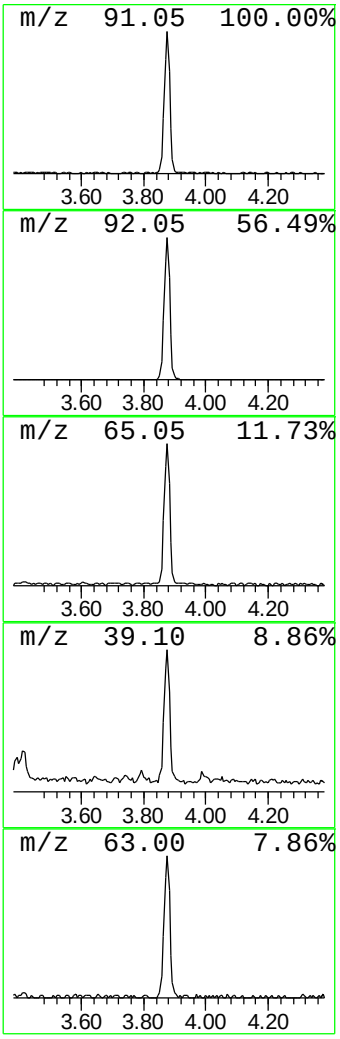
****

Peak Number 2 Toluene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.88	4.22 ng/ul	769125	1,4-Dichlorobenzene-d4	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Toluene	92	C7H8	000108-88-3	95
2	Toluene	92	C7H8	000108-88-3	94
3	Toluene	92	C7H8	000108-88-3	93
4	Toluene	92	C7H8	000108-88-3	87
5	Toluene	92	C7H8	000108-88-3	83





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Instrument : BNA_M
ClientSampleId : BFB96

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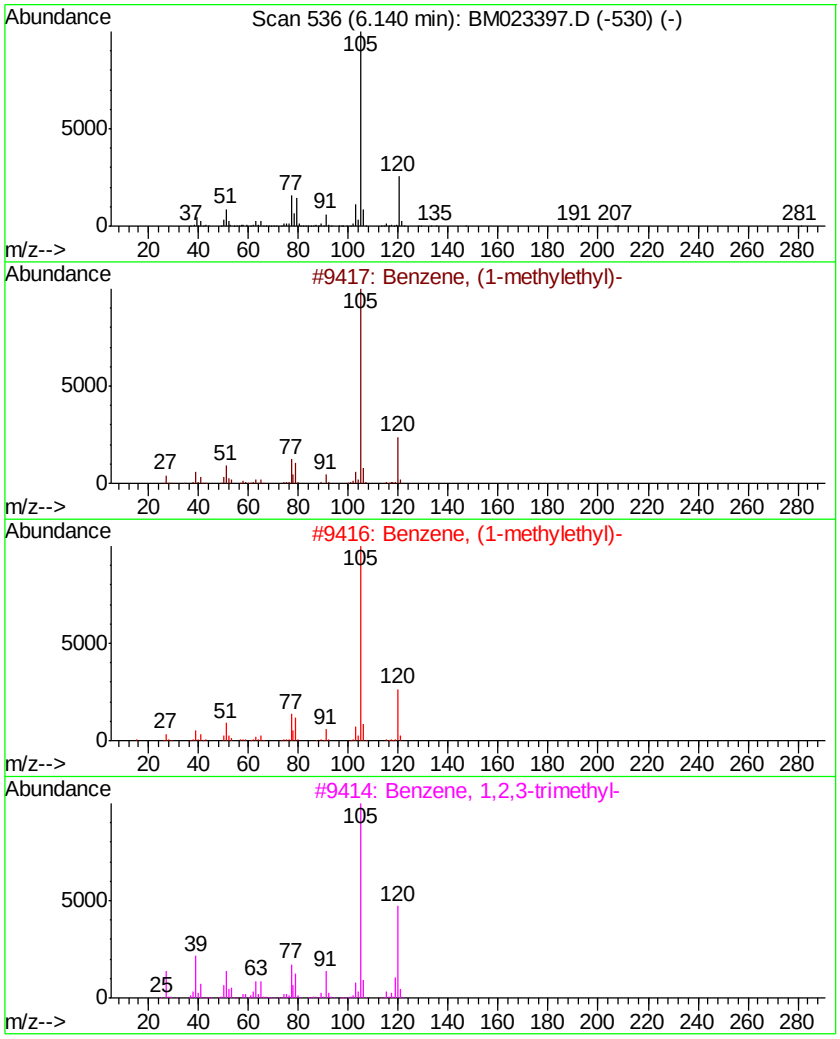
Quant Title : SVOA CALIBRATION

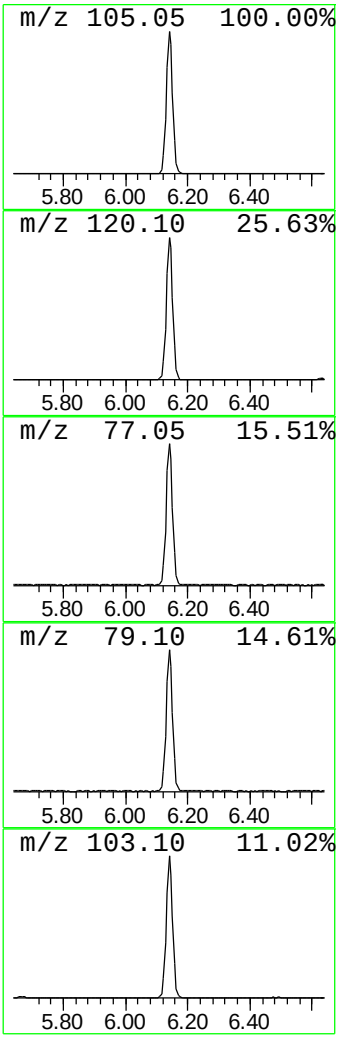
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, (1-methylethyl)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.14	14.67 ng/ul	2676300	1,4-Dichlorobenzene-d4	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	94
2	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	94
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	91
5	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	87





Library Search Compound Report

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Instrument : BNA_M
ClientSampleId : BFB96

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

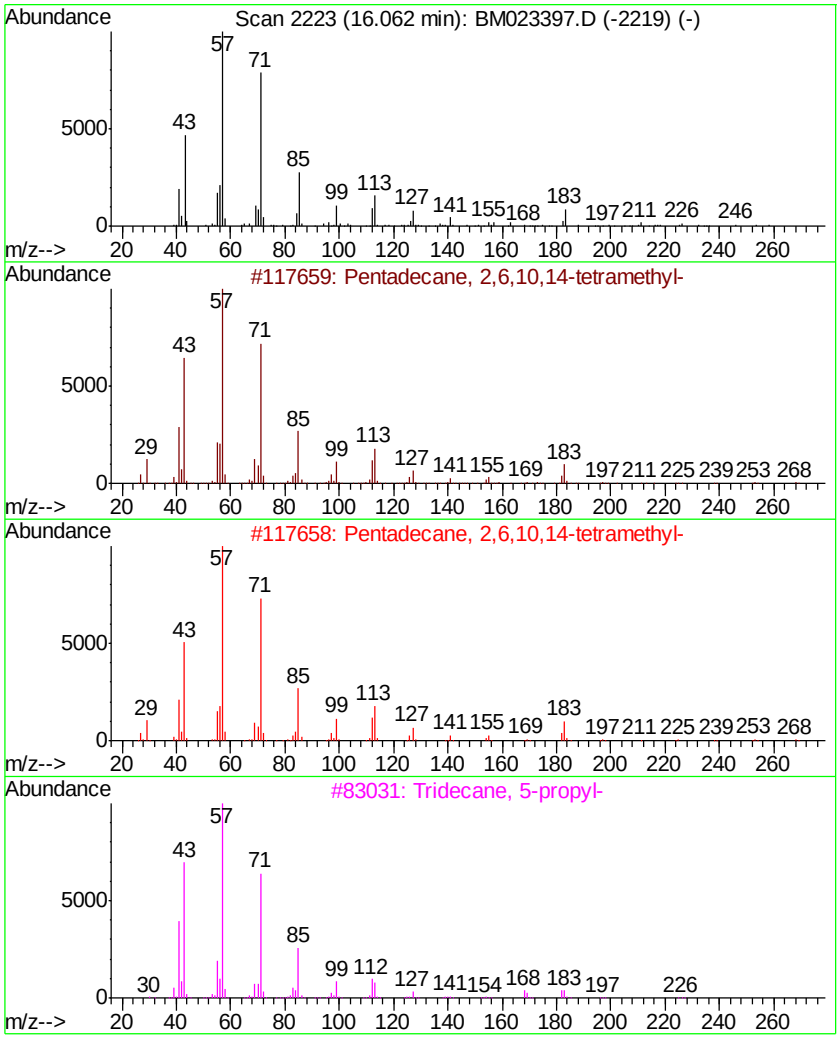
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

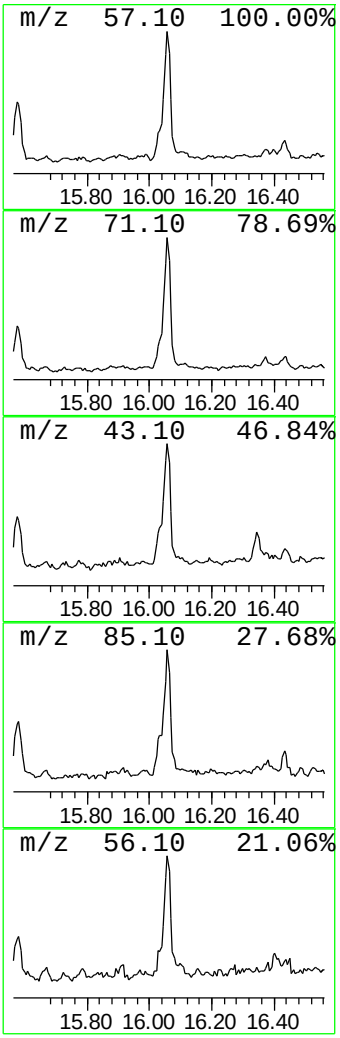
****

Peak Number 4 Pentadecane, 2,6,10,14-tetr... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.06	2.13 ng/ul	597407	Phenanthrene-d10	17.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	91
2	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	90
3	Tridecane, 5-propyl-	226	C16H34	055045-11-9	86
4	Octadecane, 2,6-dimethyl-	282	C20H42	075163-97-2	86
5	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	72





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Operator : JU
Sample : K5403-07
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFB96

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

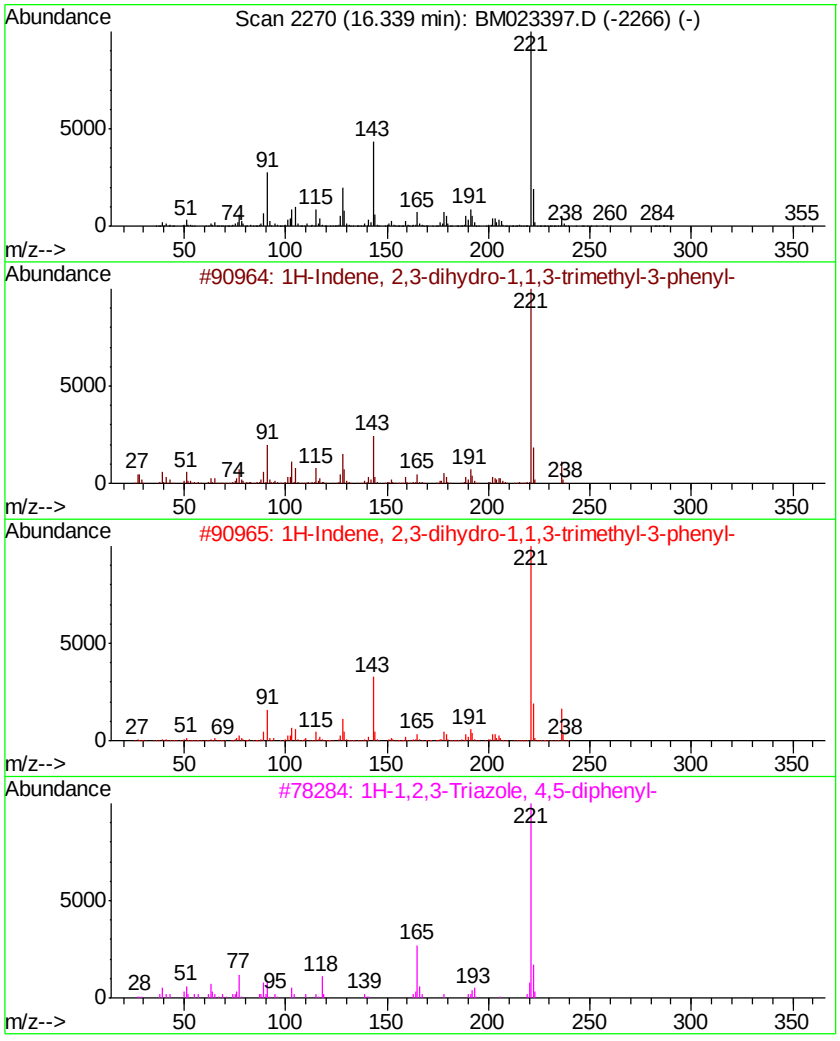
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

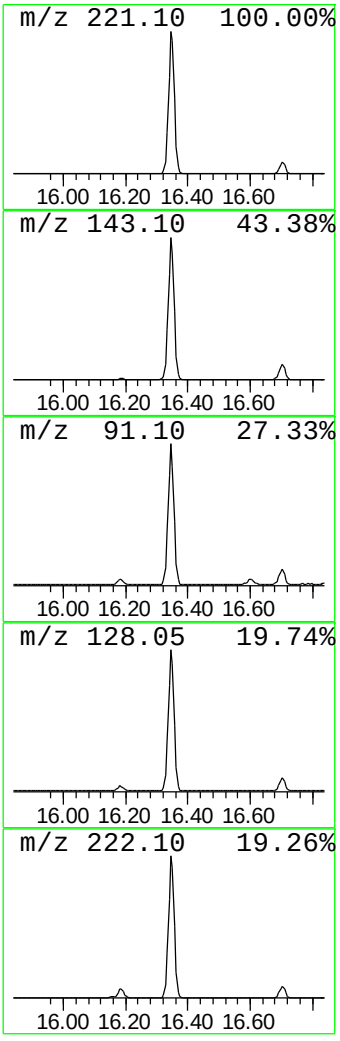
****

Peak Number 5 1H-Indene, 2,3-dihydro-1,1,... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.34	39.48 ng/ul	11084400	Phenanthrene-d10	17.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,1,3-tri...	236	C18H20	003910-35-8	93
2	1H-Indene, 2,3-dihydro-1,1,3-tri...	236	C18H20	003910-35-8	78
3	1H-1,2,3-Triazole, 4,5-diphenyl-	221	C14H11N3	005533-73-3	43
4	Carbonic acid, monoamide, N-(2-e...	221	C13H19NO2	1000339-98-3	43
5	Benz[j]isoquinoline, 3,6,8-trime...	221	C16H15N	061171-16-2	43





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM102419\
Data File : BM023397.D
Acq On : 25 Oct 2019 06:00
Operator : JU
Sample : K5403-07
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFB96

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

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

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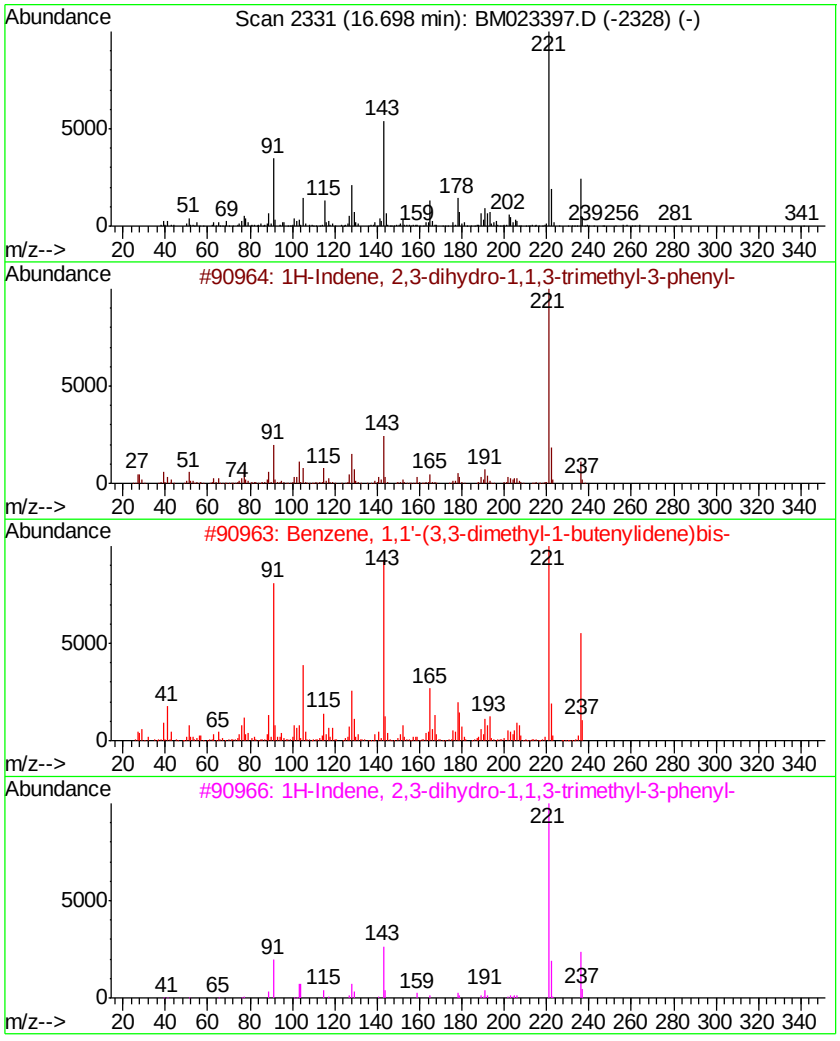
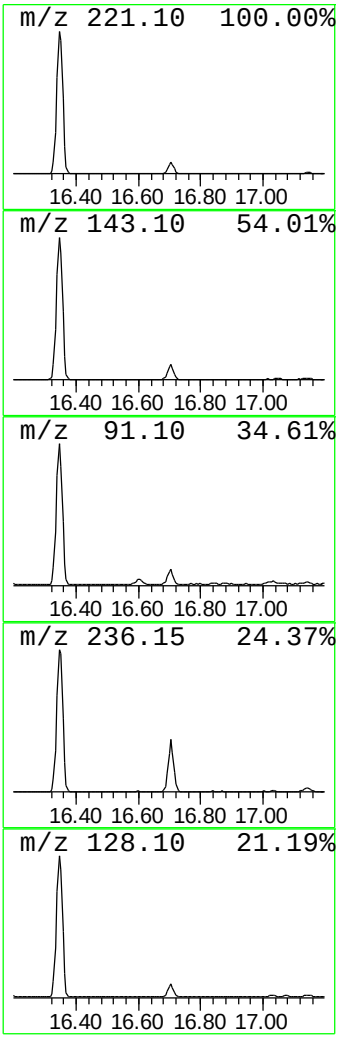
Peak Number 6 Benzene, 1,1'-(3,3-dimethyl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.70	3.80 ng/ul	1067030	Phenanthrene-d10	17.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,1,3-tri...	236	C18H20	003910-35-8	64
2	Benzene, 1,1'-(3,3-dimethyl-1-bu...	236	C18H20	023586-64-3	60
3	1H-Indene, 2,3-dihydro-1,1,3-tri...	236	C18H20	003910-35-8	58
4	1H-Indole, 1-ethyl-2-phenyl-	221	C16H15N	013228-39-2	53
5	1H-Indene, 2,3-dihydro-1,1,3-tri...	236	C18H20	003910-35-8	50





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Data File : BM023397.D
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Sample : K5403-07
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFB96

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

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

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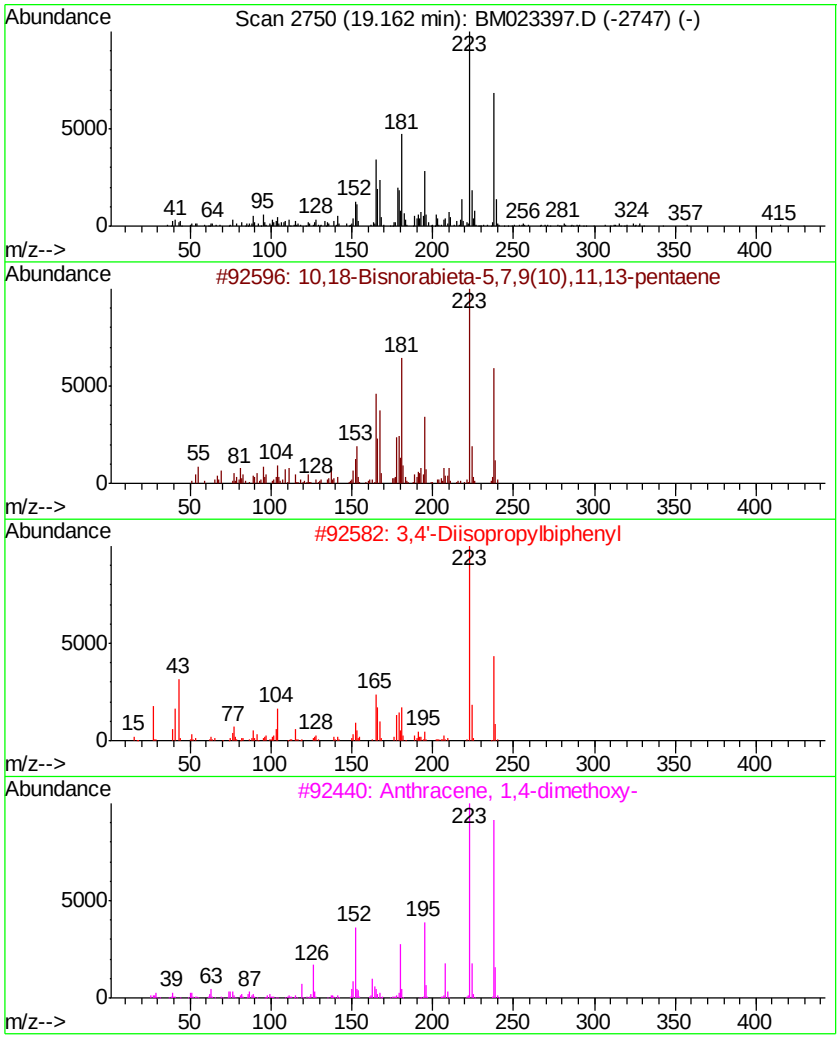
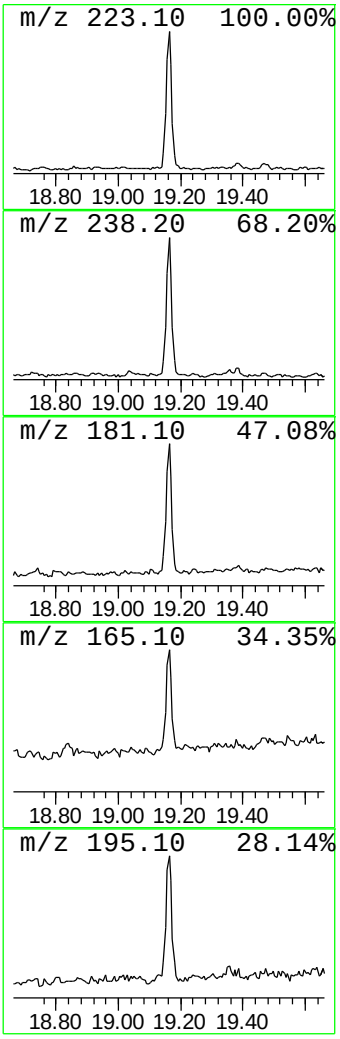
Peak Number 7 10,18-Bisnorabieta-5,7,9(10... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.16	3.33 ng/ul	701510	Chrysene-d12	21.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	10,18-Bisnorabieta-5,7,9(10),11,...	238	C18H22	006566-19-4	99
2	3,4'-Diisopropylbiphenyl	238	C18H22	061434-46-6	72
3	Anthracene, 1,4-dimethoxy-	238	C16H14O2	013076-29-4	55
4	4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	55
5	4,4'-Diisopropylbiphenyl	238	C18H22	018970-30-4	53





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Operator : JU
Sample : K5403-07
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFB96

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

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

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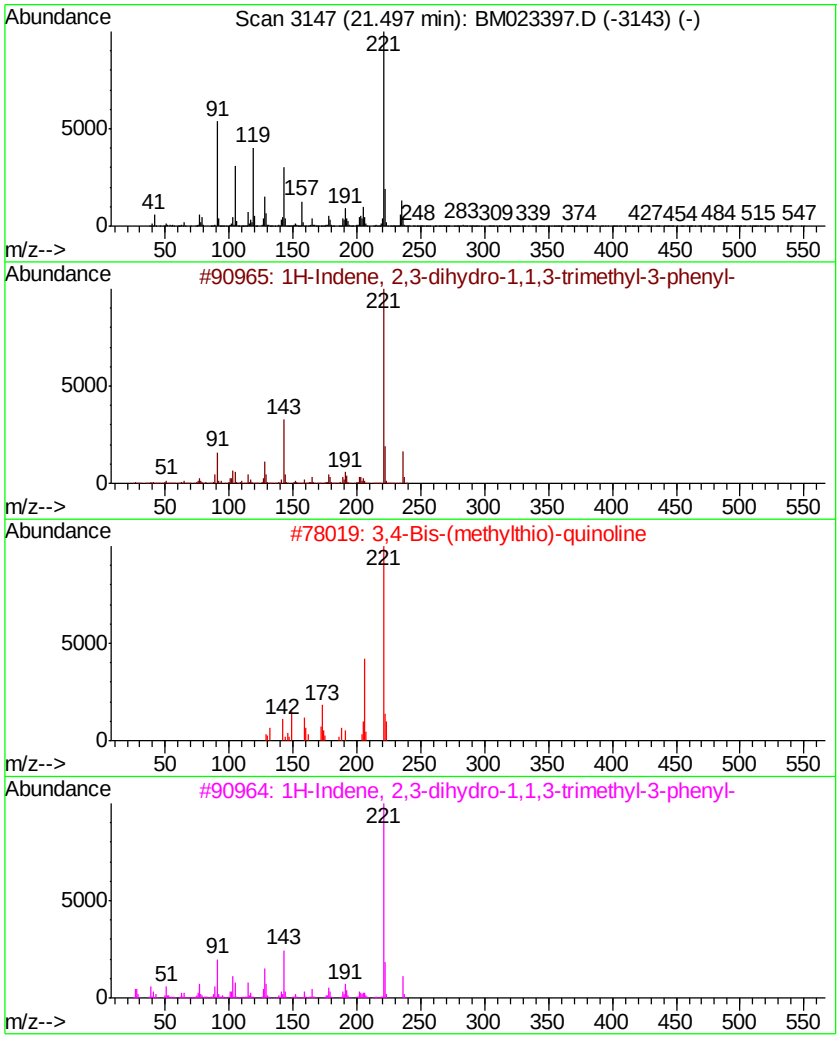
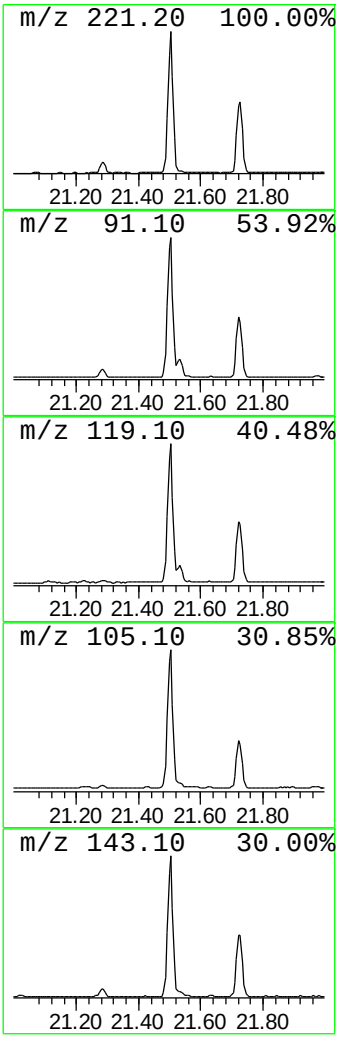
Peak Number 8 3,4-Bis-(methylthio)-quinoline Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.50	84.82 ng/ul	17870700	Chrysene-d12	21.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,1,3-tri...	236	C18H20	003910-35-8	50
2	3,4-Bis-(methylthio)-quinoline	221	C11H11NS2	074579-34-3	50
3	1H-Indene, 2,3-dihydro-1,1,3-tri...	236	C18H20	003910-35-8	47
4	3-Amino-4,6-dimethylthieno[2,3-b...	221	C10H11N3OS	067795-42-0	47
5	Boron, diethyl(5-ethyl-4,6-nonan...	250	C15H31BN2	136705-03-8	47





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Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFB96

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

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

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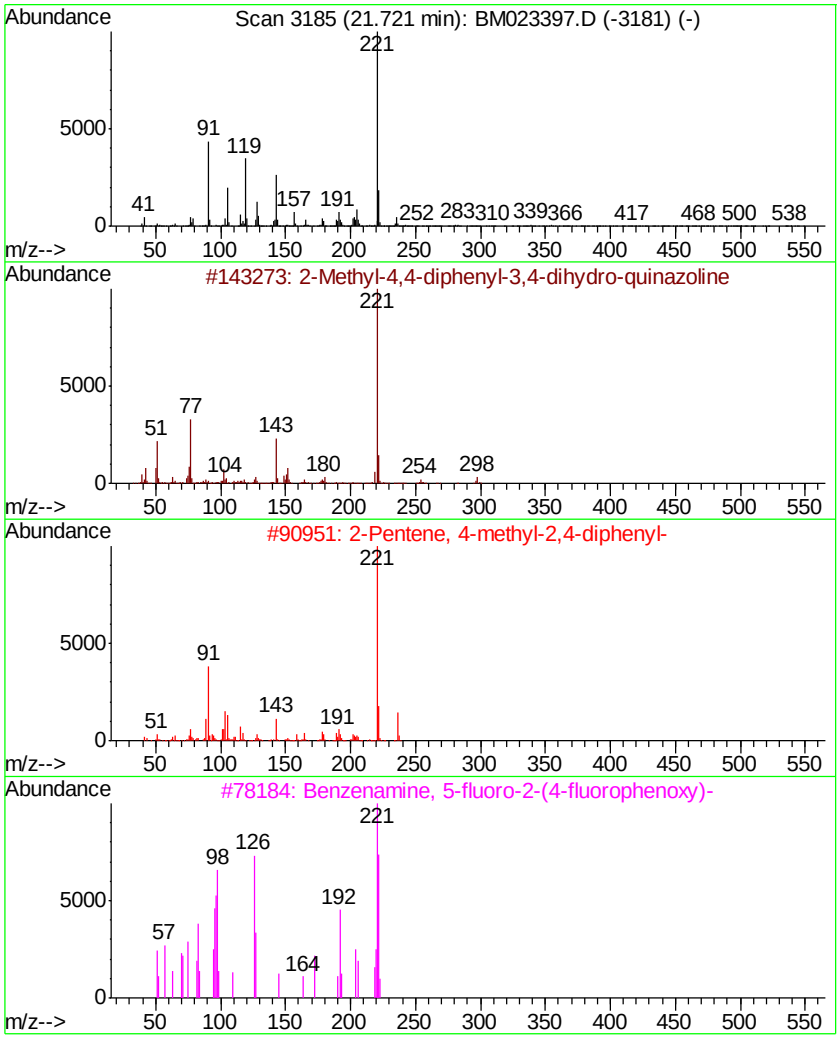
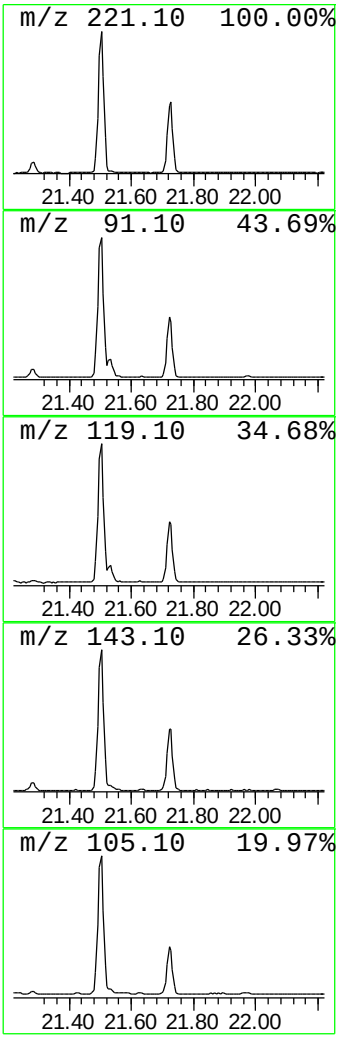
Peak Number 9 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.72	37.83 ng/ul	7970720	Chrysene-d12	21.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methyl-4,4-diphenyl-3,4-dihydr...	298	C21H18N2	1000300-40-0	47
2	2-Pentene, 4-methyl-2,4-diphenyl-	236	C18H20	006258-73-7	40
3	Benzenamine, 5-fluoro-2-(4-fluor...	221	C12H9F2NO	020653-64-9	38
4	Boron, diethyl(5-ethyl-4,6-nonan...	250	C15H31BN2	136705-03-8	38
5	4-Hydroxy-beta-phenylcinnamonitrile	221	C15H11NO	016281-90-6	32





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Acq On : 25 Oct 2019 06:00
Operator : JU
Sample : K5403-07
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFB96

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

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

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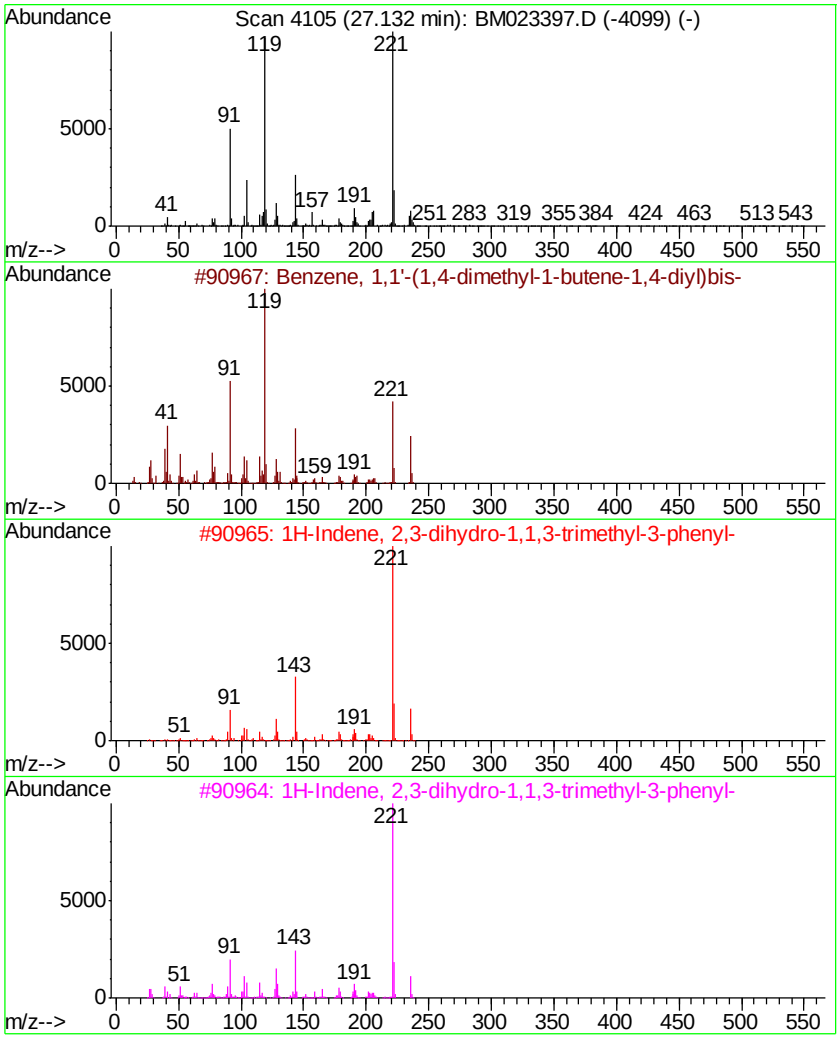
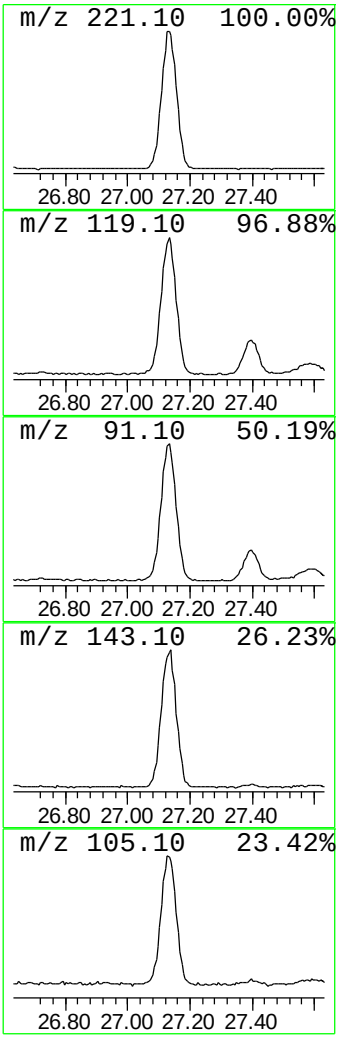
Peak Number 10 Benzene, 1,1'-(1,4-dimethyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.13	23.08 ng/ul	6959030	Perylene-d12	23.43

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,1'-(1,4-dimethyl-1-bu...	236	C18H20	052161-54-3	50
2	1H-Indene, 2,3-dihydro-1,1,3-tri...	236	C18H20	003910-35-8	49
3	1H-Indene, 2,3-dihydro-1,1,3-tri...	236	C18H20	003910-35-8	43
4	1H-Indene, 2,3-dihydro-1,1,3-tri...	236	C18H20	003910-35-8	37
5	2-Methyl-4,4-diphenyl-3,4-dihydr...	298	C21H18N2	1000300-40-0	35





Tentatively Identified Compound (LSC) summary

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Data File : BM023397.D
Acq On : 25 Oct 2019 06:00
Operator : JU
Sample : K5403-07
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BFB96

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM102319MA.M
Quant Title : SVOA CALIBRATION

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(DEL) Alkane: Str...	3.09	7.4	ng/ul	1342720	1	7.63	3648530	20.0
Toluene	3.88	4.2	ng/ul	769125	1	7.63	3648530	20.0
Benzene, (1-methy...	6.14	14.7	ng/ul	2676300	1	7.63	3648530	20.0
Pentadecane, 2,6,...	16.06	2.1	ng/ul	597407	4	17.03	5615140	20.0
1H-Indene, 2,3-di...	16.34	39.5	ng/ul	11084400	4	17.03	5615140	20.0
Benzene, 1,1'-(3,...	16.70	3.8	ng/ul	1067030	4	17.03	5615140	20.0
10,18-Bisnorabiet...	19.16	3.3	ng/ul	701510	5	21.23	4213600	20.0
3,4-Bis-(methylth...	21.50	84.8	ng/ul	17870700	5	21.23	4213600	20.0
unknown-01	21.72	37.8	ng/ul	7970720	5	21.23	4213600	20.0
Benzene, 1,1'-(1,...	27.13	23.1	ng/ul	6959030	6	23.43	6029870	20.0