LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112719\
Data File : BM023963.D
Acq On : 28 Nov 2019 21:25
Operator : JU
Sample : K5854-17
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AD3

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-BM112719MA.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	3.034	22	25	39	rVB	110089	209085	1.70%	0.143%
2	3.657	125	131	135	rVB	71488	113285	0.92%	0.078%
3	3.740	140	145	157	rVB	685083	942606	7.68%	0.647%
4	4.646	294	299	304	rBV	63974	100342	0.82%	0.069%
5	5.028	356	364	370	rVV	265787	391882	3.19%	0.269%
6	5.157	380	386	398	rBV	626478	1325544	10.81%	0.910%
7	5.522	439	448	455	rBV2	350873	653759	5.33%	0.449%
8	5.993	524	528	534	rVB2	29887	49404	0.40%	0.034%
9	6.475	604	610	618	rVB2	98823	183657	1.50%	0.126%
10	6.587	618	629	634	rBV	341254	535377	4.36%	0.367%
11	6.675	634	644	662	rVV2	1694873	3555841	28.99%	2.440%
12	6.834	666	671	677	rVV	1412607	2219931	18.10%	1.523%
13	6.881	677	679	690	rVB	134320	205901	1.68%	0.141%
14	7.022	696	703	714	rVB	1938919	3056901	24.92%	2.098%
15	7.140	714	723	732	rVB	470879	792675	6.46%	0.544%
16	7.487	774	782	791	rBV	1867089	2940038	23.97%	2.018%
17	7.569	791	796	798	rVV2	50675	87171	0.71%	0.060%
18	7.598	798	801	812	rVB2	106213	185999	1.52%	0.128%
19	7.857	839	845	850	rVB	73204	119183	0.97%	0.082%
20	7.981	861	866	870	rBV	42577	61813	0.50%	0.042%
21	8.034	870	875	881	rVB2	121078	191621	1.56%	0.131%
22	8.128	885	891	895	rBV2	158661	255784	2.09%	0.176%
23	8.204	898	904	911	rVV	2037630	3250964	26.50%	2.231%
24	8.263	911	914	916	rVV4	76452	118956	0.97%	0.082%
25	8.310	918	922	929	rVB	211199	361625	2.95%	0.248%
26	8.434	939	943	948	rBV	49451	75318	0.61%	0.052%
27	8.486	948	952	959	rVB	59437	94155	0.77%	0.065%
28	8.586	961	969	972	rBV	158212	251968	2.05%	0.173%
29	8.639	972	978	987	rVV	1678116	2844485	23.19%	1.952%
30	8.722	987	992	999	rVV	97622	158810	1.29%	0.109%
31	8.828	1007	1010	1018	rVB6	26033	54326	0.44%	0.037%
32	8.910	1018	1024	1029	rBV2	28253	51442	0.42%	0.035%
33	9.022	1039	1043	1049	rBV3	34987	73919	0.60%	0.051%
34	9.163	1063	1067	1075	rVB	70053	123702	1.01%	0.085%

LSC Area Percent Report

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Sample : K5854-17
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AD3

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-BM112719MA.M
Title : SVOA CALIBRATION

35	9.351	1091	1099	1107	rBV	1409148	2416350	19.70%	1.658%
36	9.516	1123	1127	1131	rVV	40291	59675	0.49%	0.041%
37	9.551	1131	1133	1142	rVB8	43800	68819	0.56%	0.047%
38	9.633	1142	1147	1149	rBV2	51793	81998	0.67%	0.056%
39	9.663	1149	1152	1157	rVB2	62865	103344	0.84%	0.071%
40	9.716	1157	1161	1163	rBV3	41887	55916	0.46%	0.038%
41	9.892	1182	1191	1200	rVV	2284347	3912399	31.90%	2.685%
42	9.975	1201	1205	1213	rVB	55173	99393	0.81%	0.068%
43	10.257	1245	1253	1259	rVV	2723590	4459211	36.35%	3.060%
44	10.310	1259	1262	1271	rVB	201346	325067	2.65%	0.223%
45	10.404	1271	1278	1289	rBV	1715434	3023813	24.65%	2.075%
46	10.828	1343	1350	1355	rBV8	24118	54084	0.44%	0.037%
47	11.016	1378	1382	1391	rBV6	27220	68145	0.56%	0.047%
48	11.116	1394	1399	1405	rVB6	37367	71268	0.58%	0.049%
49	11.175	1405	1409	1416	rBV8	25218	61627	0.50%	0.042%
50	11.304	1426	1431	1439	rBV5	62566	155561	1.27%	0.107%
51	11.457	1449	1457	1460	rBV3	56030	94264	0.77%	0.065%
52	11.663	1484	1492	1497	rBV9	38309	104640	0.85%	0.072%
53	11.716	1497	1501	1510	rVV	117632	206221	1.68%	0.142%
54	11.816	1512	1518	1521	rVV6	28767	54391	0.44%	0.037%
55	11.857	1521	1525	1530	rVV2	56209	92478	0.75%	0.063%
56	11.933	1530	1538	1547	rVB	213337	413412	3.37%	0.284%
57	12.157	1571	1576	1580	rBV	91468	170923	1.39%	0.117%
58	12.198	1580	1583	1588	rVB3	55733	82750	0.67%	0.057%
59	12.557	1640	1644	1648	rBV7	35569	63062	0.51%	0.043%
60	12.686	1662	1666	1670	rBV2	76453	110823	0.90%	0.076%
61	12.986	1708	1717	1724	rBV3	184580	354953	2.89%	0.244%
62	13.080	1729	1733	1737	rVB4	38094	60800	0.50%	0.042%
63	13.304	1764	1771	1773	rBV5	99698	185937	1.52%	0.128%
64	13.463	1794	1798	1803	rVV3	69771	121418	0.99%	0.083%
65	13.516	1803	1807	1810	rBV4	61444	92022	0.75%	0.063%
66	13.569	1810	1816	1820	rVV	4208714	5617062	45.79%	3.855%
67	13.616	1820	1824	1828	rVV	983370	1346570	10.98%	0.924%
68	13.657	1828	1831	1834	rVB2	92351	126689	1.03%	0.087%
69	13.833	1854	1861	1873	rBV	5099727	7741235	63.11%	5.312%
70	14.074	1897	1902	1906	rBV	191194	245072	2.00%	0.168%
71	14.145	1906	1914	1920	rBV	4136965	6182693	50.41%	4.243%

LSC Area Percent Report

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Sample : K5854-17
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AD3

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

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Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Title : SVOA CALIBRATION

72	14.380	1948	1954	1968	rBV	1129752	2027714	16.53%	1.391%
73	14.598	1986	1991	1995	rBV2	342659	558603	4.55%	0.383%
74	14.698	2005	2008	2014	rVB4	136322	209375	1.71%	0.144%
75	14.768	2017	2020	2024	rBV5	78976	144176	1.18%	0.099%
76	14.939	2046	2049	2053	rBV4	74130	106349	0.87%	0.073%
77	15.033	2061	2065	2069	rVB	224286	256784	2.09%	0.176%
78	15.151	2079	2085	2090	rBV	6411887	9188494	74.91%	6.305%
79	15.292	2104	2109	2114	rBV	923073	1273885	10.39%	0.874%
80	15.439	2132	2134	2139	rVB2	195600	254993	2.08%	0.175%
81	15.898	2208	2212	2214	rBV2	248649	331433	2.70%	0.227%
82	15.921	2214	2216	2224	rVB2	321472	439190	3.58%	0.301%
83	16.421	2298	2301	2305	rBV	191576	263840	2.15%	0.181%
84	16.686	2342	2346	2350	rBV2	698756	909987	7.42%	0.624%
85	16.745	2353	2356	2360	rVV	284554	363180	2.96%	0.249%
86	16.792	2360	2364	2367	rVB	176583	210632	1.72%	0.145%
87	16.904	2377	2383	2387	rBV	4490026	6738640	54.94%	4.624%
88	16.939	2387	2389	2394	rVV	569657	736994	6.01%	0.506%
89	16.998	2394	2399	2409	rVV	6869354	9805058	79.94%	6.728%
90	17.209	2431	2435	2441	rVB	522129	680383	5.55%	0.467%
91	17.609	2499	2503	2512	rVB	448156	672425	5.48%	0.461%
92	17.827	2535	2540	2546	rBV	1186250	1671988	13.63%	1.147%
93	18.968	2731	2734	2742	rVB	970800	1212854	9.89%	0.832%
94	19.115	2757	2759	2763	rBV2	409167	497107	4.05%	0.341%
95	19.309	2786	2792	2800	rBV2	7795736	12265983	100.00%	8.417%
96	20.850	3051	3054	3058	rVB	1279061	1344357	10.96%	0.923%
97	21.044	3084	3087	3092	rBV	1624683	1993981	16.26%	1.368%
98	21.109	3093	3098	3102	rBV	6018982	7855526	64.04%	5.391%
99	23.121	3435	3440	3445	rBV	6619131	10846144	88.42%	7.443%
100	23.256	3458	3463	3474	rVB	4749050	8744552	71.29%	6.001%

Sum of corrected areas: 145726181

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112719\
Data File : BM023963.D
Acq On : 28 Nov 2019 21:25
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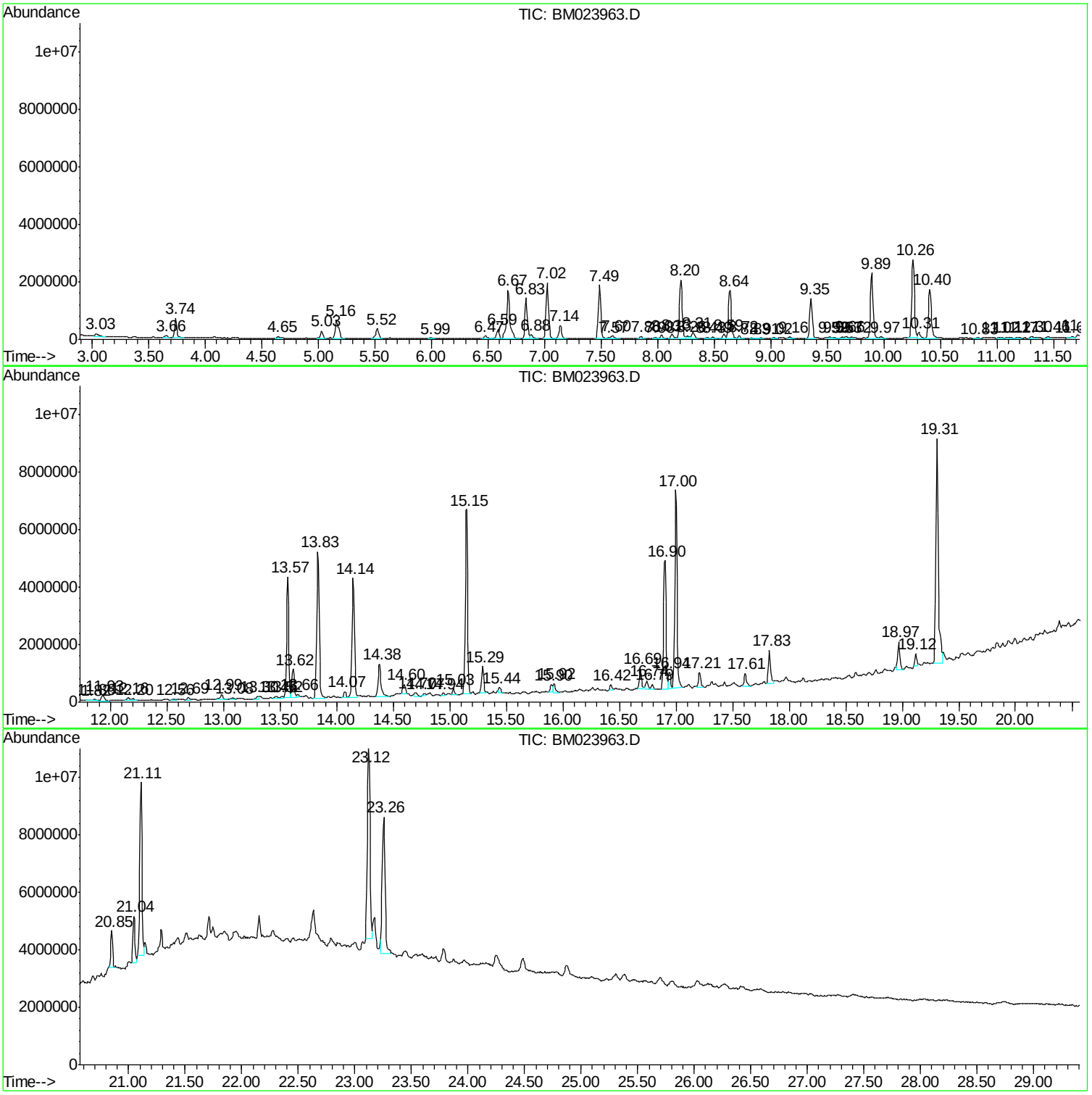
Sample : K5854-17
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0AD3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.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_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

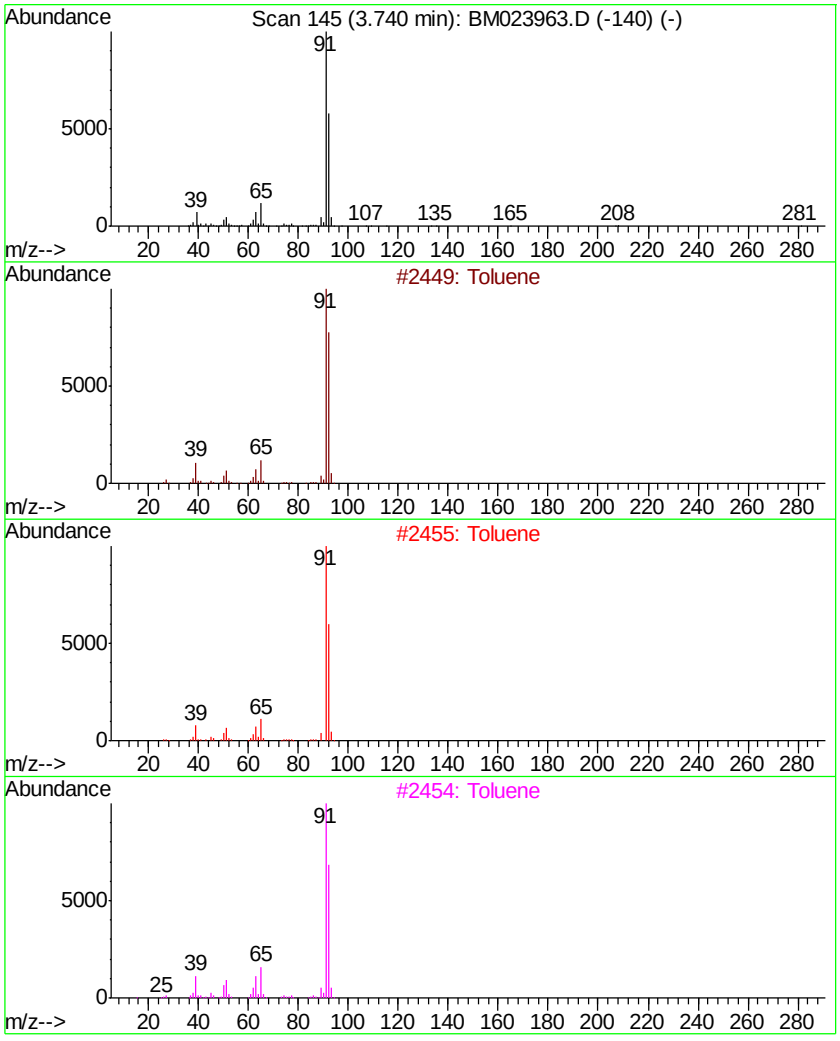
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

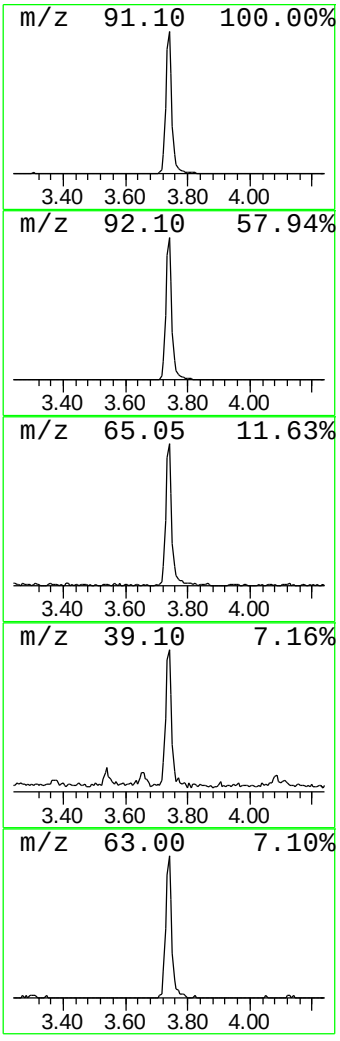
****

Peak Number 1 Toluene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.74	6.41 ng/ul	942606	1,4-Dichlorobenzene-d4	7.49

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	91
4	Toluene	92	C7H8	000108-88-3	91
5	Spiro[2,4]hepta-4,6-diene	92	C7H8	000765-46-8	87





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112719\
Data File : BM023963.D
Acq On : 28 Nov 2019 21:25
Operator : JU
Sample : K5854-17
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AD3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

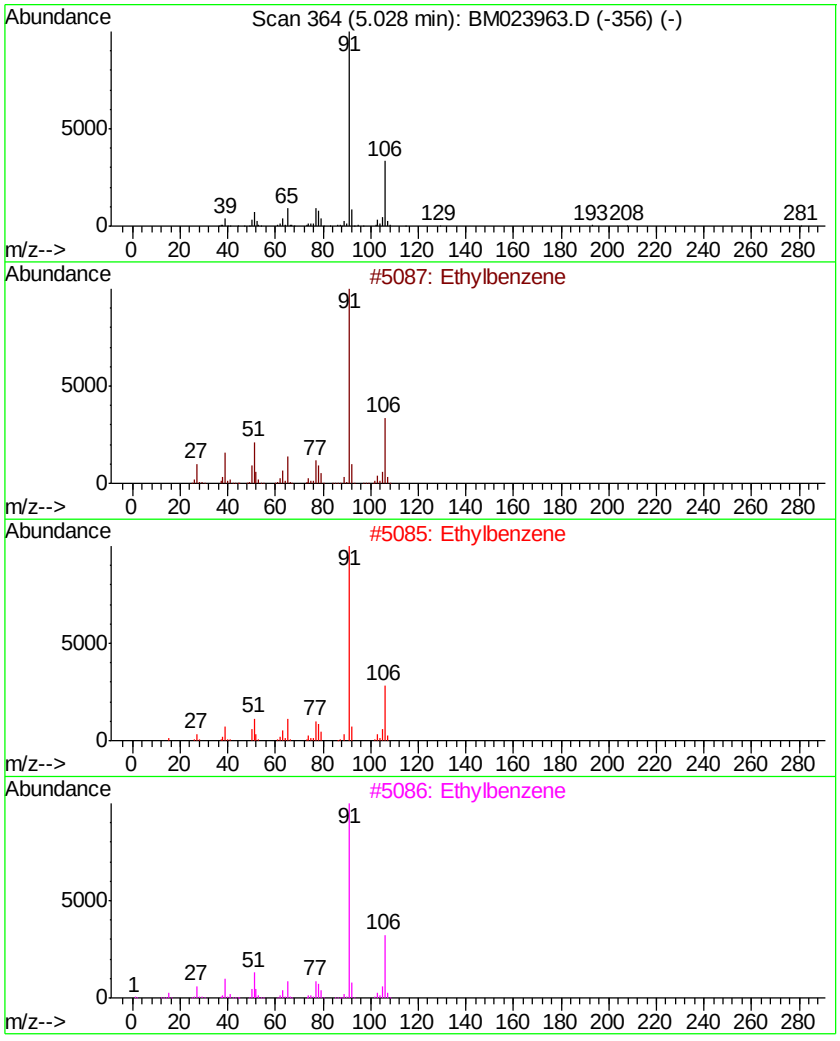
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

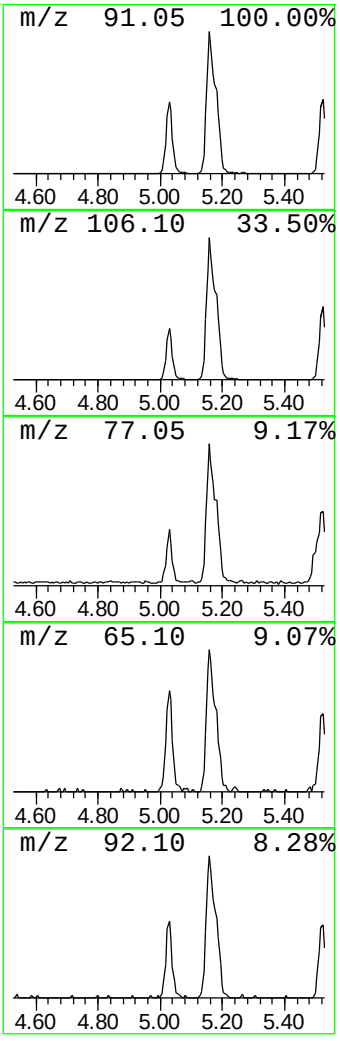
****

Peak Number 2 Ethylbenzene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.03	2.67 ng/ul	391882	1,4-Dichlorobenzene-d4	7.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethylbenzene	106	C8H10	000100-41-4	91
2	Ethylbenzene	106	C8H10	000100-41-4	90
3	Ethylbenzene	106	C8H10	000100-41-4	90
4	Ethylbenzene	106	C8H10	000100-41-4	90
5	p-Xylene	106	C8H10	000106-42-3	68





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112719\
Data File : BM023963.D
Acq On : 28 Nov 2019 21:25
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ALS Vial : 47 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AD3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

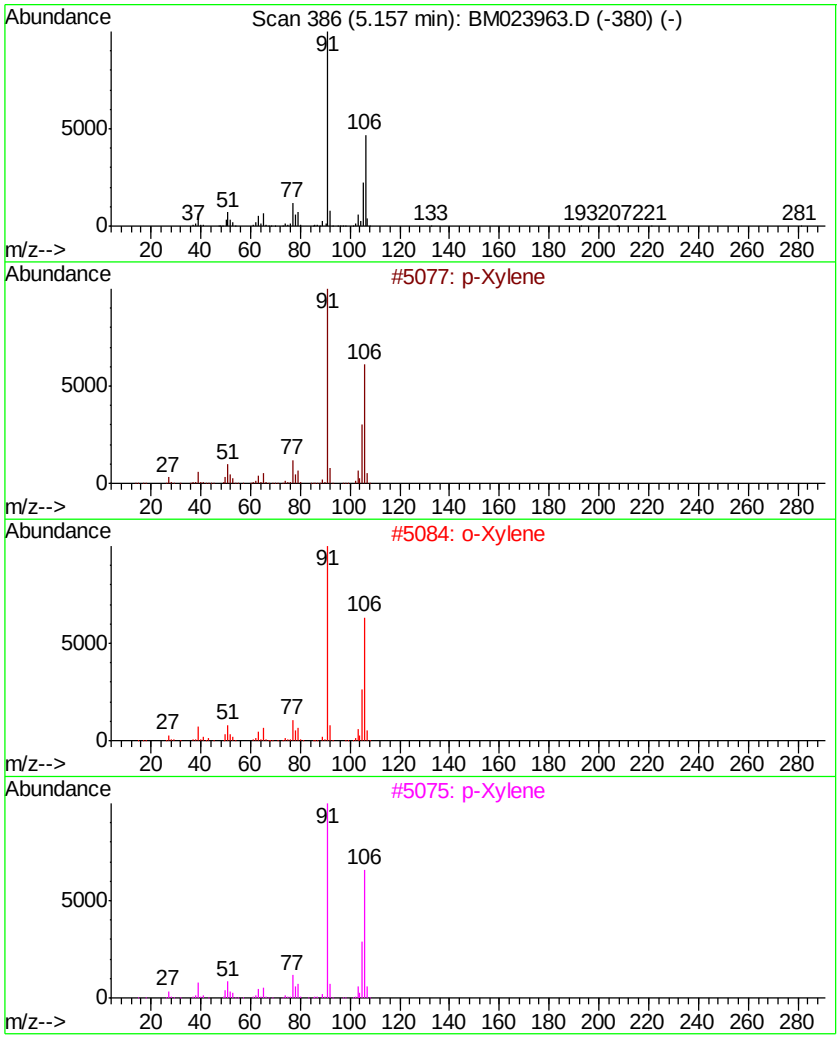
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

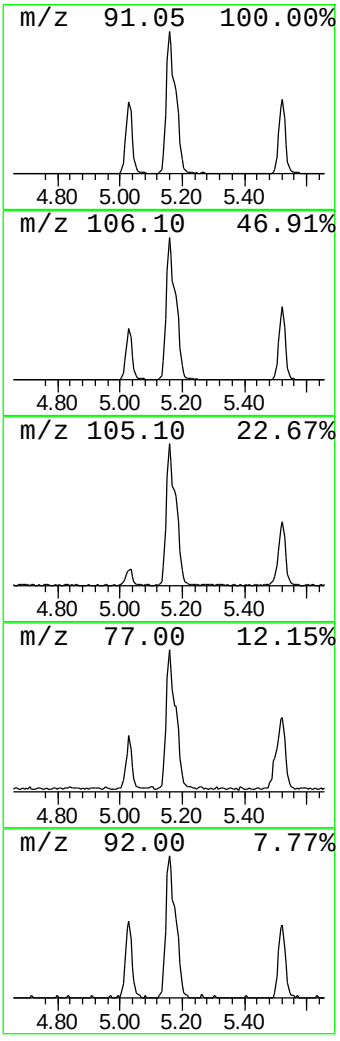
****

Peak Number 3 p-Xylene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.16	9.02 ng/ul	1325540	1,4-Dichlorobenzene-d4	7.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Xylene	106	C8H10	000106-42-3	97
2	o-Xylene	106	C8H10	000095-47-6	95
3	p-Xylene	106	C8H10	000106-42-3	95
4	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	95
5	o-Xylene	106	C8H10	000095-47-6	94





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Instrument : BNA_M
ClientSampleId : C0AD3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

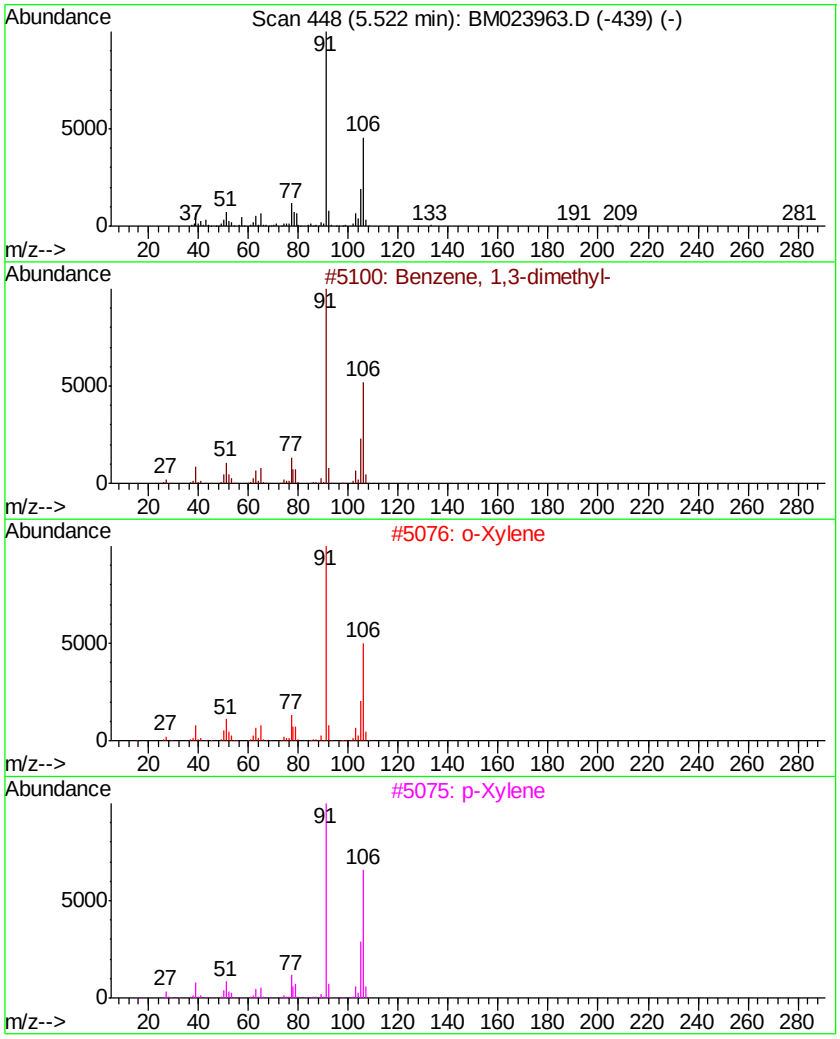
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

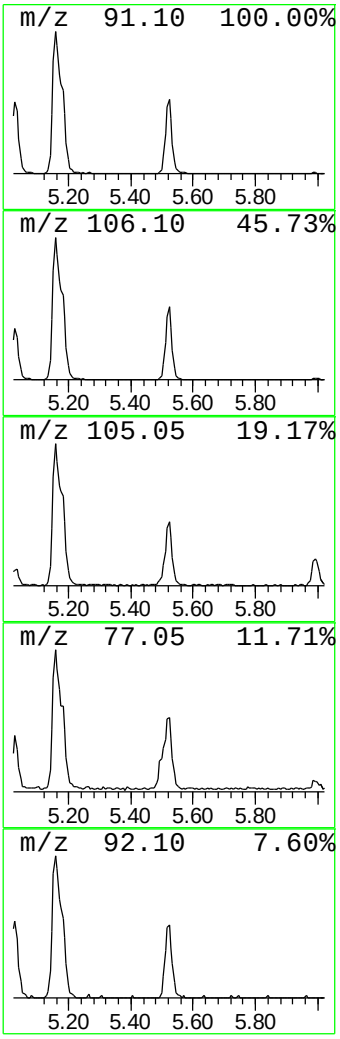
****

Peak Number 4 Benzene, 1,3-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.52	4.45 ng/ul	653759	1,4-Dichlorobenzene-d4	7.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	95
2	o-Xylene	106	C8H10	000095-47-6	95
3	p-Xylene	106	C8H10	000106-42-3	95
4	p-Xylene	106	C8H10	000106-42-3	94
5	p-Xylene	106	C8H10	000106-42-3	93





Library Search Compound Report

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Data File : BM023963.D
Acq On : 28 Nov 2019 21:25
Operator : JU
Sample : K5854-17
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AD3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

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

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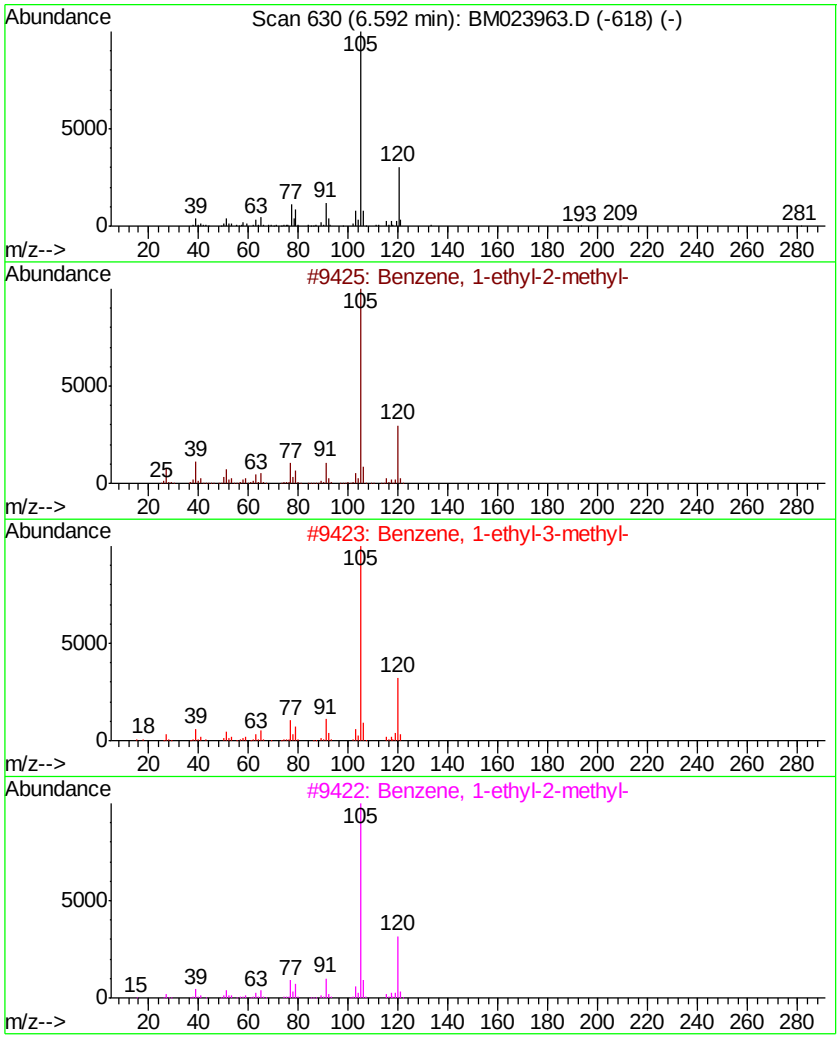
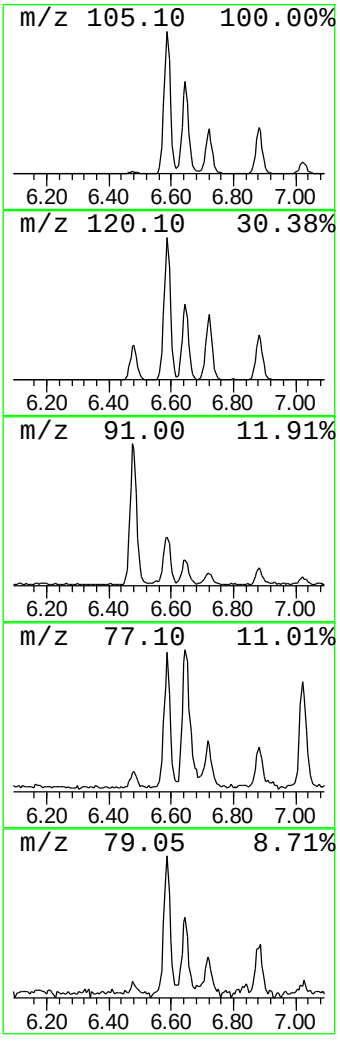
Peak Number 5 Benzene, 1-ethyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.59	3.64 ng/ul	535377	1,4-Dichlorobenzene-d4	7.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM112719\
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ALS Vial : 47 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AD3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

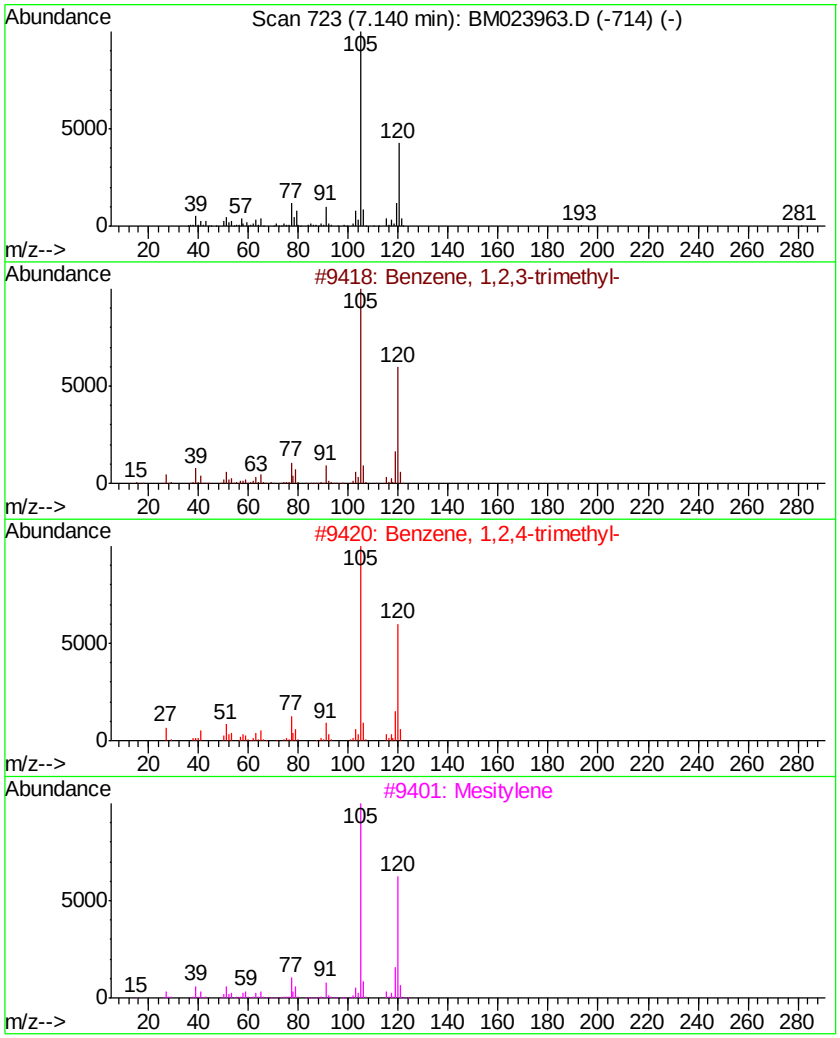
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

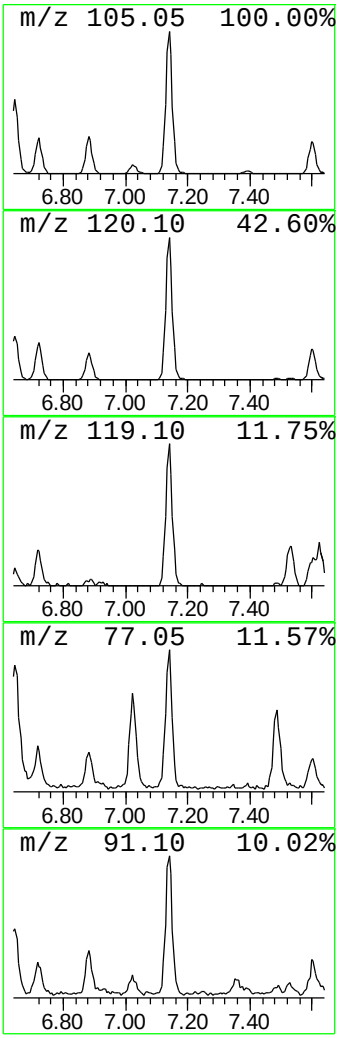
****

Peak Number 6 Benzene, 1,2,3-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.14	5.39 ng/ul	792675	1,4-Dichlorobenzene-d4	7.49

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
3	Mesitylene	120	C9H12	000108-67-8	95
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94





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Data File : BM023963.D
Acq On : 28 Nov 2019 21:25
Operator : JU
Sample : K5854-17
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AD3

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

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

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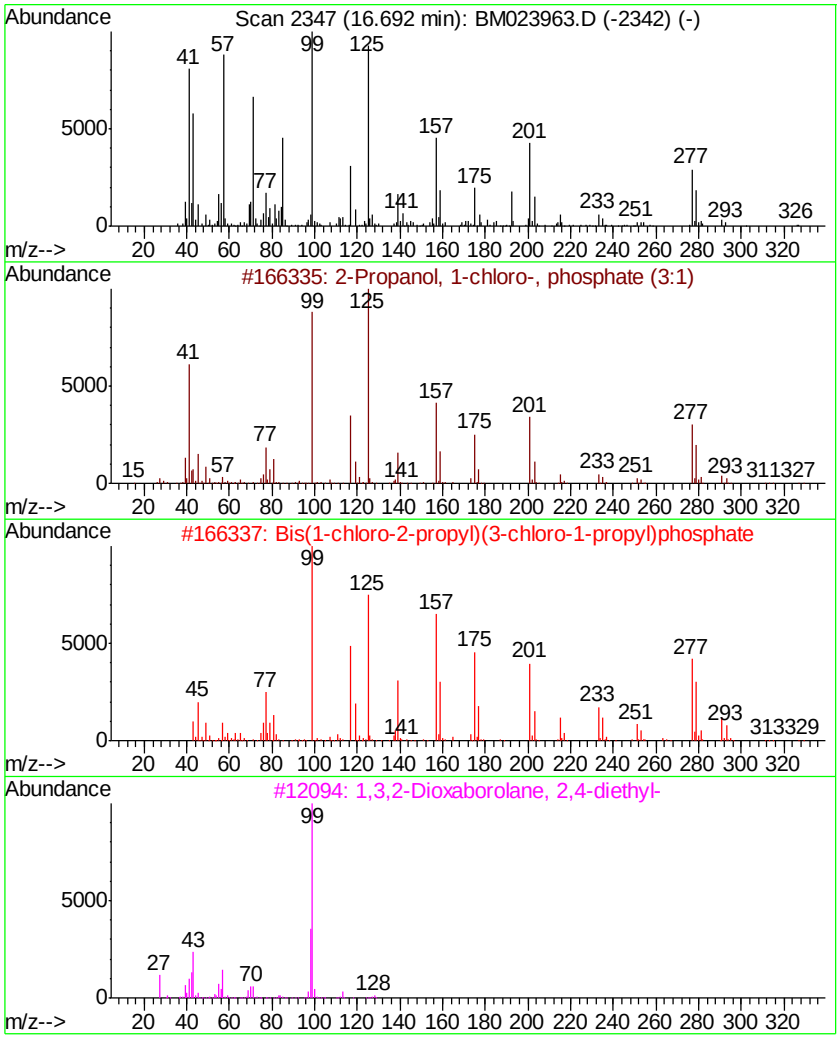
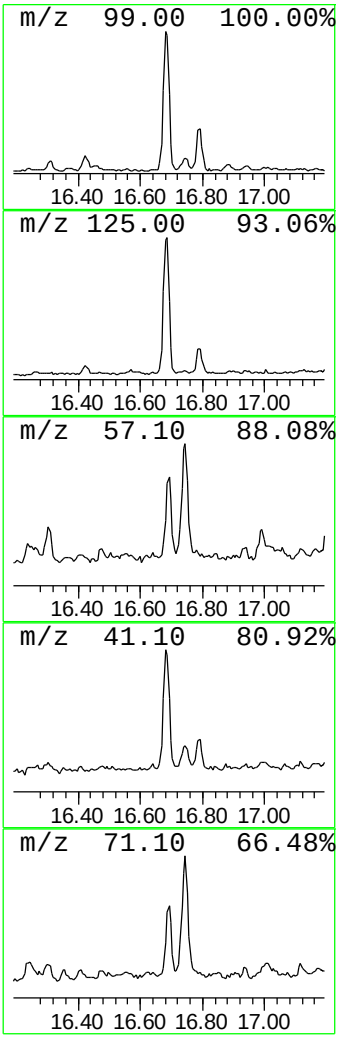
Peak Number 7 2-Propanol, 1-chloro-, phos... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.69	2.70 ng/ul	909987	Phenanthrene-d10	16.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propanol, 1-chloro-, phosphate...	326	C9H18Cl3O4P	013674-84-5	55
2	Bis(1-chloro-2-propyl)(3-chloro-...	326	C9H18Cl3O4P	137909-40-1	22
3	1,3,2-Dioxaborolane, 2,4-diethyl-	128	C6H13BO2	057633-63-3	11
4	Isopropylphosphonic acid, dicylc...	260	C13H25O3P	1000273-18-4	10
5	2-Amino-4-hydroxy-6-methylpyrimi...	125	C5H7N3O	003977-29-5	10





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Operator : JU
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ALS Vial : 47 Sample Multiplier: 1

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ClientSampleId : C0AD3

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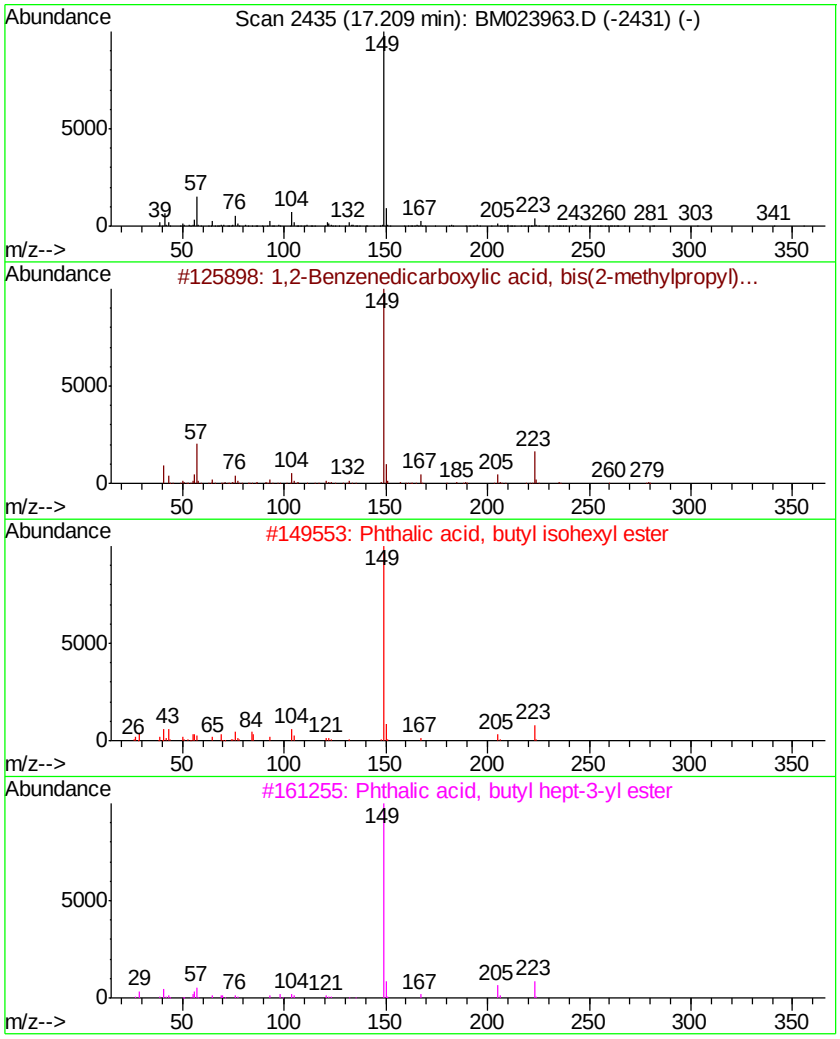
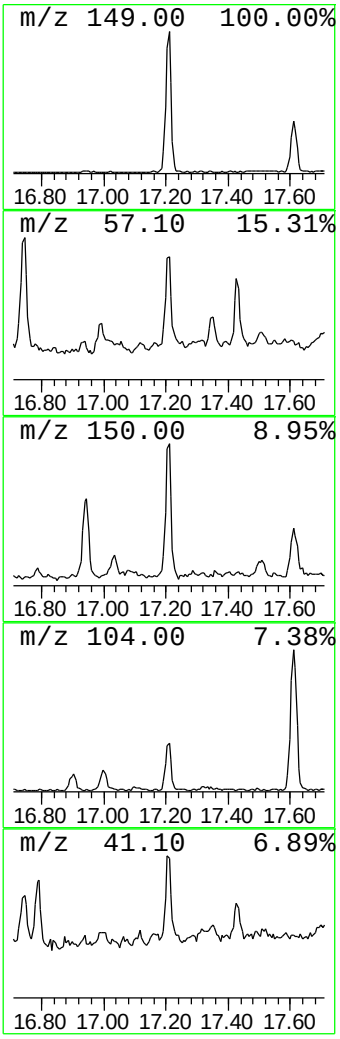
Peak Number 8 1,2-Benzenedicarboxylic aci... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.21	2.02 ng/ul	680383	Phenanthrene-d10	16.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Benzenedicarboxylic acid, bi...	278	C16H22O4	000084-69-5	83
2	Phthalic acid, butyl isohexyl ester	306	C18H26O4	1000309-03-6	78
3	Phthalic acid, butyl hept-3-yl e...	320	C19H28O4	1000356-98-7	78
4	1,2-Benzenedicarboxylic acid, bi...	278	C16H22O4	000084-69-5	78
5	Phthalic acid, butyl hexyl ester	306	C18H26O4	1000308-99-5	78





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Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AD3

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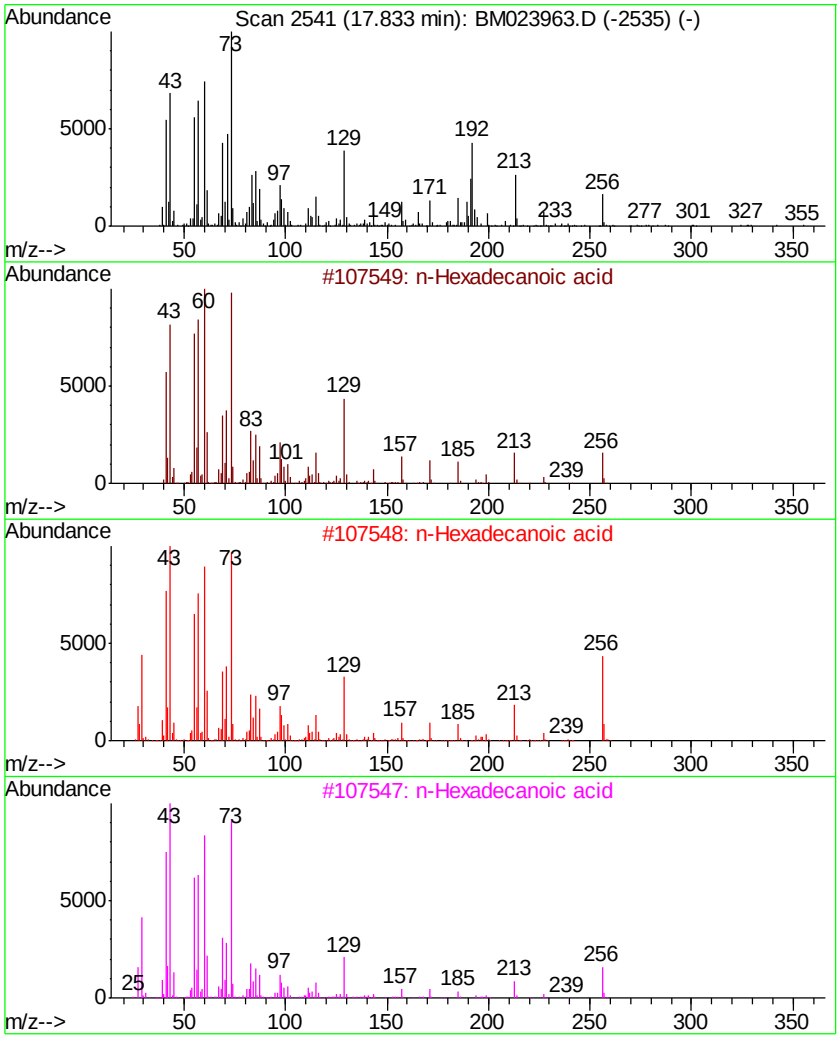
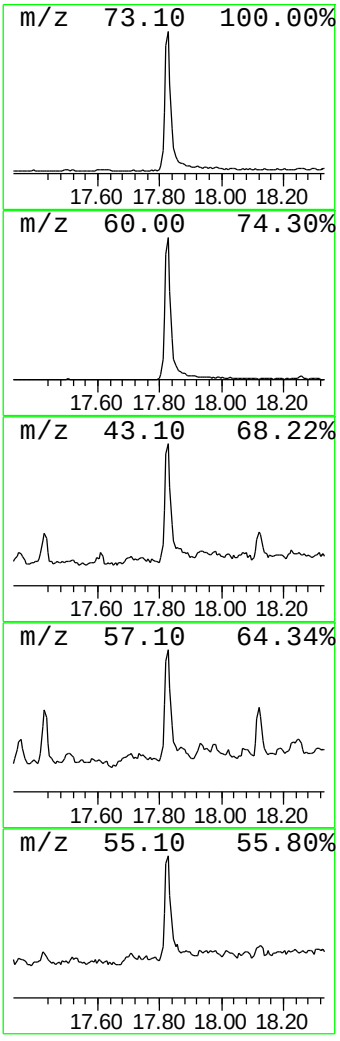
Peak Number 9 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.83	4.96 ng/ul	1671990	Phenanthrene-d10	16.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	68
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	64





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Operator : JU
Sample : K5854-17
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C0AD3

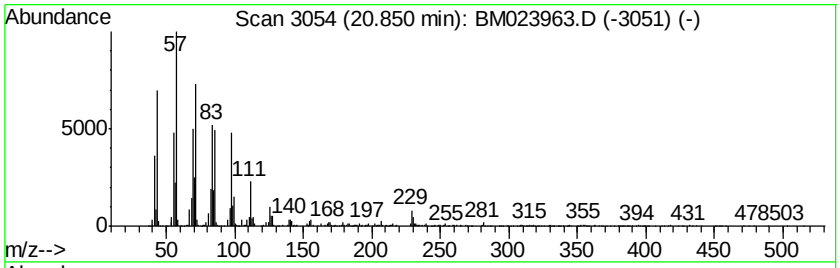
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM112719MA.M
Quant Title : SVOA CALIBRATION

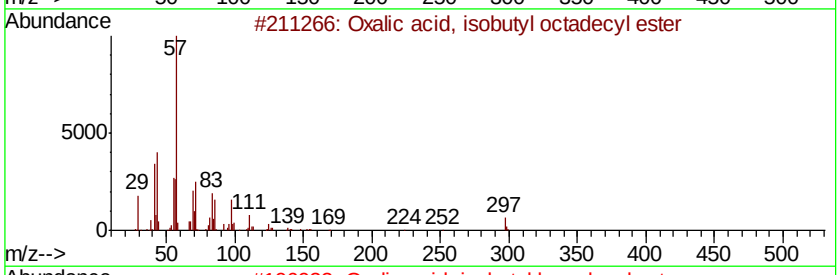
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

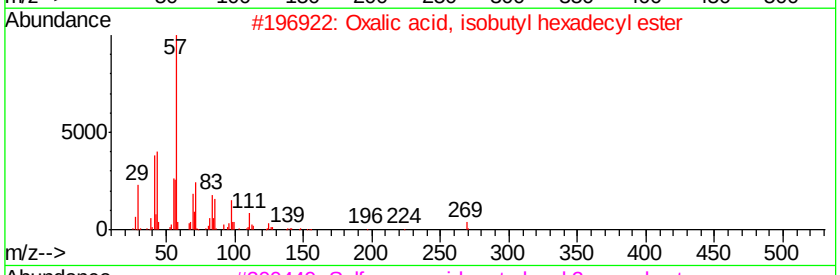
****
Peak Number 10 Oxalic acid, isobutyl octad... Concentration Rank 7

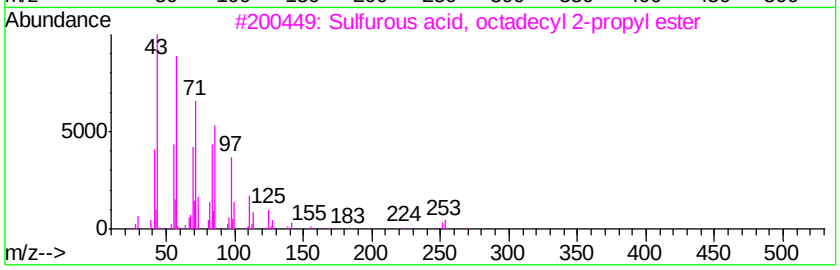
R.T.	EstConc	Area	Relative to ISTD	R.T.
20.85	3.42 ng/ul	1344360	Chrysene-d12	21.11

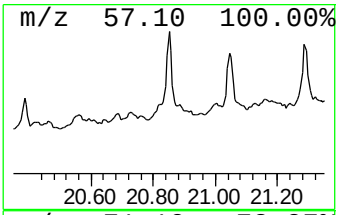
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxalic acid, isobutyl octadecyl ...	398	C24H46O4	1000309-38-3	91
2	Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	91
3	Sulfurous acid, octadecyl 2-prop...	376	C21H44O3S	1000309-12-7	91
4	Oxalic acid, isobutyl heptadecyl...	384	C23H44O4	1000309-38-2	90
5	Ethanol, 2-(hexadecyloxy)-	286	C18H38O2	002136-71-2	87

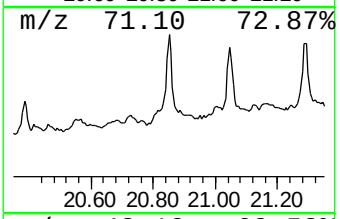


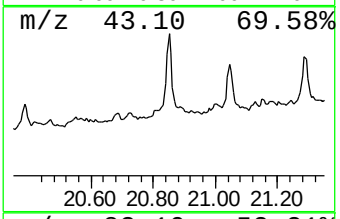


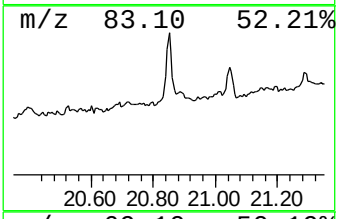


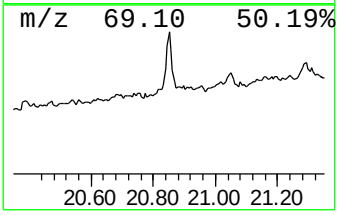












Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM112719\
Data File : BM023963.D
Acq On : 28 Nov 2019 21:25
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ALS Vial : 47 Sample Multiplier: 1

Instrument : BNA_M
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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
Toluene	3.74	6.4	ng/ul	942606	1	7.49	2940040	20.0
Ethylbenzene	5.03	2.7	ng/ul	391882	1	7.49	2940040	20.0
p-Xylene	5.16	9.0	ng/ul	1325540	1	7.49	2940040	20.0
Benzene, 1,3-dime...	5.52	4.5	ng/ul	653759	1	7.49	2940040	20.0
Benzene, 1-ethyl-...	6.59	3.6	ng/ul	535377	1	7.49	2940040	20.0
Benzene, 1,2,3-tr...	7.14	5.4	ng/ul	792675	1	7.49	2940040	20.0
2-Propanol, 1-chl...	16.69	2.7	ng/ul	909987	4	16.90	6738640	20.0
1,2-Benzenedicarb...	17.21	2.0	ng/ul	680383	4	16.90	6738640	20.0
n-Hexadecanoic acid	17.83	5.0	ng/ul	1671990	4	16.90	6738640	20.0
Oxalic acid, isob...	20.85	3.4	ng/ul	1344360	5	21.11	7855530	20.0