LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX102220\
Data File : VX019033.D
Acq On : 22 Oct 2020 17:30
Operator : JC/SP
Sample : L4417-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-14

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102120W.M
Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.776	108	113	120	rVB	65906	88520	1.82%	0.212%
2	2.380	205	212	217	rBV	31285	61333	1.26%	0.147%
3	2.825	278	285	289	rBV	42978	94714	1.95%	0.227%
4	2.873	289	293	300	rVB	37533	72959	1.50%	0.175%
5	3.154	329	339	351	rVB	39981	92284	1.90%	0.221%
6	4.373	531	539	551	rVB2	67607	191135	3.93%	0.459%
7	5.464	702	718	727	rBV	168120	487006	10.00%	1.168%
8	5.629	736	745	772	rVB	271643	850891	17.48%	2.041%
9	6.031	801	811	826	rBV	172962	448708	9.22%	1.076%
10	6.324	849	859	870	rVB2	91811	280939	5.77%	0.674%
11	6.830	932	942	952	rBV	429344	1037864	21.32%	2.490%
12	6.909	952	955	965	rVB3	21715	53566	1.10%	0.129%
13	7.239	1001	1009	1016	rVB2	27319	61750	1.27%	0.148%
14	7.440	1031	1042	1052	rVB6	34129	99966	2.05%	0.240%
15	7.556	1052	1061	1066	rBV	29706	59576	1.22%	0.143%
16	7.934	1112	1123	1131	rBV7	22911	87389	1.80%	0.210%
17	8.037	1131	1140	1151	rVB	146756	298397	6.13%	0.716%
18	8.159	1151	1160	1169	rBV2	248603	540959	11.11%	1.298%
19	8.238	1169	1173	1184	rVV4	38473	81875	1.68%	0.196%
20	8.354	1185	1192	1198	rVV3	24089	51579	1.06%	0.124%
21	8.549	1216	1224	1233	rVB4	47000	105374	2.16%	0.253%
22	8.696	1233	1248	1258	rBV	997208	1675674	34.42%	4.020%
23	9.202	1327	1331	1336	rBV	44476	67959	1.40%	0.163%
24	9.549	1383	1388	1394	rVB3	24565	53718	1.10%	0.129%
25	10.098	1473	1478	1483	rBV	915331	1251071	25.70%	3.001%
26	10.232	1496	1500	1509	rVB	331174	449106	9.23%	1.077%
27	10.342	1515	1518	1524	rVB	53091	77405	1.59%	0.186%
28	11.000	1621	1626	1633	rBV	3147860	4004277	82.25%	9.606%
29	11.122	1641	1646	1651	rBV	837377	1045526	21.48%	2.508%
30	11.341	1677	1682	1691	rVB	3941542	4868338	100.00%	11.679%
31	11.439	1693	1698	1702	rVV	314392	382656	7.86%	0.918%
32	11.488	1702	1706	1714	rVB	378854	475427	9.77%	1.141%
33	11.652	1727	1733	1742	rBV	99374	136781	2.81%	0.328%
34	11.793	1751	1756	1761	rVB	279864	339390	6.97%	0.814%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX102220\
Data File : VX019033.D
Acq On : 22 Oct 2020 17:30
Operator : JC/SP
Sample : L4417-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-14

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102120W.M
Title : SW846 8260

35	11.902	1768	1774	1776	rBV	256065	325156	6.68%	0.780%
36	11.927	1776	1778	1784	rVB	292964	359766	7.39%	0.863%
37	12.061	1794	1800	1806	rBV	1193875	1424156	29.25%	3.417%
38	12.128	1806	1811	1816	rBV	599491	738814	15.18%	1.772%
39	12.286	1831	1837	1843	rBV	2442788	3166891	65.05%	7.597%
40	12.353	1843	1848	1858	rVB2	555156	1052859	21.63%	2.526%
41	12.445	1858	1863	1867	rBV	174171	208243	4.28%	0.500%
42	12.500	1867	1872	1879	rVB	310625	374620	7.70%	0.899%
43	12.567	1879	1883	1890	rBV	47562	66423	1.36%	0.159%
44	12.652	1890	1897	1900	rBV	1808103	2116272	43.47%	5.077%
45	12.683	1900	1902	1907	rVV	285833	311007	6.39%	0.746%
46	12.738	1907	1911	1919	rVB	697306	943790	19.39%	2.264%
47	12.884	1929	1935	1940	rVB2	79198	117136	2.41%	0.281%
48	12.963	1940	1948	1952	rBV	1431253	1841667	37.83%	4.418%
49	13.000	1952	1954	1960	rVB	152407	169596	3.48%	0.407%
50	13.061	1960	1964	1970	rBV2	98103	139347	2.86%	0.334%
51	13.152	1973	1979	1982	rBV2	136286	199639	4.10%	0.479%
52	13.207	1985	1988	1998	rVB	608740	719991	14.79%	1.727%
53	13.298	1998	2003	2004	rBV	58474	68502	1.41%	0.164%
54	13.329	2004	2008	2014	rVB2	1604194	2135117	43.86%	5.122%
55	13.408	2019	2021	2026	rVV	71526	83078	1.71%	0.199%
56	13.463	2026	2030	2036	rVB	128474	162984	3.35%	0.391%
57	13.542	2038	2043	2046	rBV	66998	92238	1.89%	0.221%
58	13.585	2046	2050	2052	rVV	155539	212241	4.36%	0.509%
59	13.622	2052	2056	2060	rVV3	250975	396061	8.14%	0.950%
60	13.682	2060	2066	2067	rVV2	88846	153620	3.16%	0.369%
61	13.701	2067	2069	2076	rVB2	200415	288185	5.92%	0.691%
62	13.817	2080	2088	2096	rBV	450463	576729	11.85%	1.384%
63	13.890	2096	2100	2105	rVB	63297	82682	1.70%	0.198%
64	13.963	2105	2112	2116	rBV2	135483	194512	4.00%	0.467%
65	14.012	2116	2120	2124	rVV	87273	100870	2.07%	0.242%
66	14.115	2132	2137	2141	rBV	189778	230583	4.74%	0.553%
67	14.256	2155	2160	2169	rBV	161065	285832	5.87%	0.686%
68	14.359	2173	2177	2182	rVV	129055	163983	3.37%	0.393%
69	14.414	2182	2186	2190	rVB	139678	176454	3.62%	0.423%
70	14.518	2198	2203	2208	rBV2	96869	143684	2.95%	0.345%
71	14.676	2217	2229	2233	rBV	263838	413824	8.50%	0.993%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX102220\
Data File : VX019033.D
Acq On : 22 Oct 2020 17:30
Operator : JC/SP
Sample : L4417-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-14

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102120W.M
Title : SW846 8260

72	14.816	2247	2252	2257	rBV	539868	709551	14.57%	1.702%
73	15.225	2313	2319	2322	rBV	53276	89608	1.84%	0.215%
74	15.450	2353	2356	2361	rVB	43615	62378	1.28%	0.150%
75	15.524	2361	2368	2373	rBV	77971	127878	2.63%	0.307%
76	15.664	2382	2391	2394	rBV	94621	153955	3.16%	0.369%
77	15.700	2394	2397	2404	rVB	52370	78832	1.62%	0.189%
78	15.865	2417	2424	2426	rBV2	34448	61753	1.27%	0.148%
79	15.896	2426	2429	2443	rVB2	45614	86763	1.78%	0.208%
80	16.042	2447	2453	2464	rVB	58895	102654	2.11%	0.246%

Sum of corrected areas: 41684036

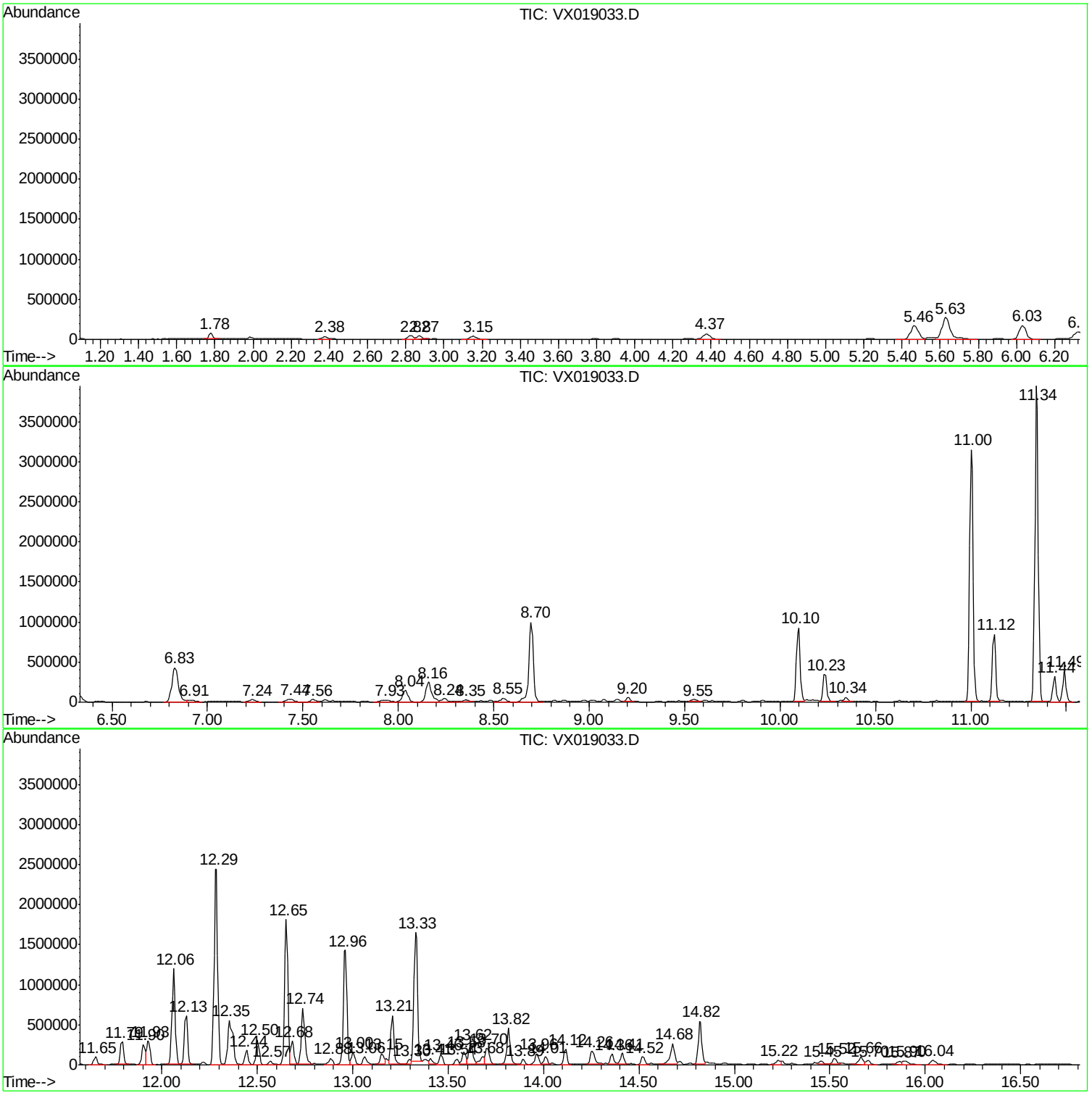
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX102220\
Data File : VX019033.D
Acq On : 22 Oct 2020 17:30
Operator : JC/SP
Sample : L4417-10
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-14

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102120W.M
Quant Title : SW846 8260

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



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX102220\
Data File : VX019033.D
Acq On : 22 Oct 2020 17:30
Operator : JC/SP
Sample : L4417-10
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-14

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102120W.M
Quant Title : SW846 8260

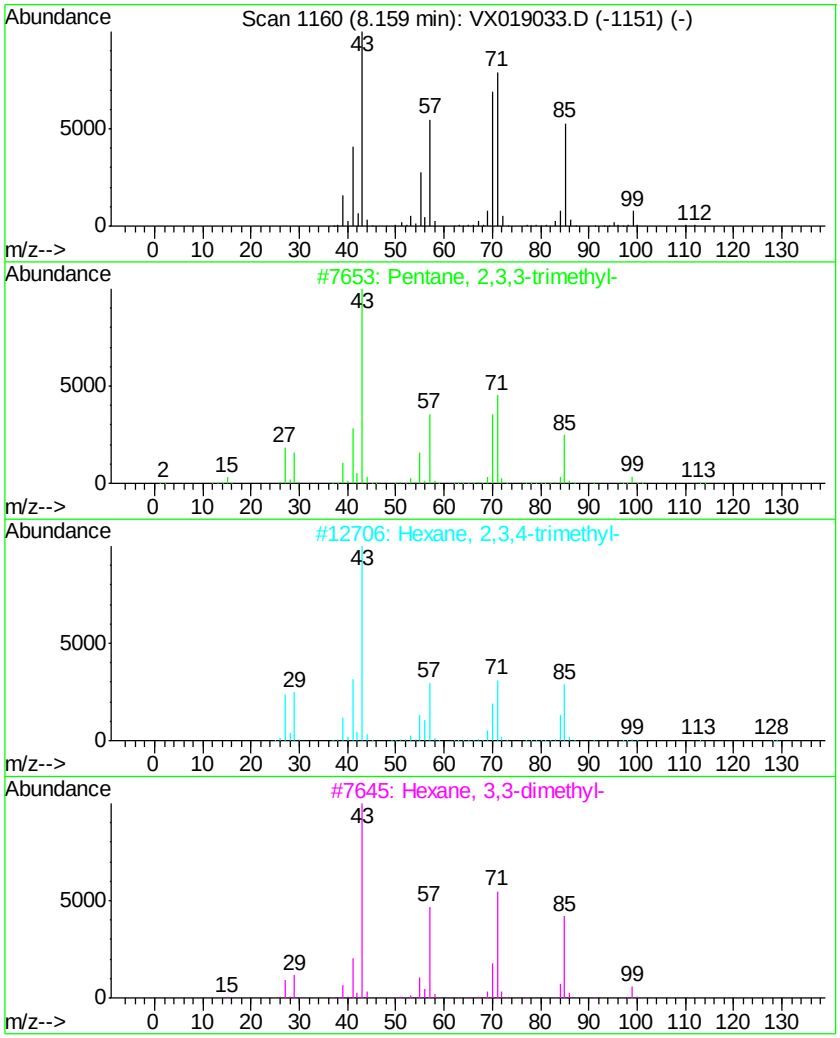
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

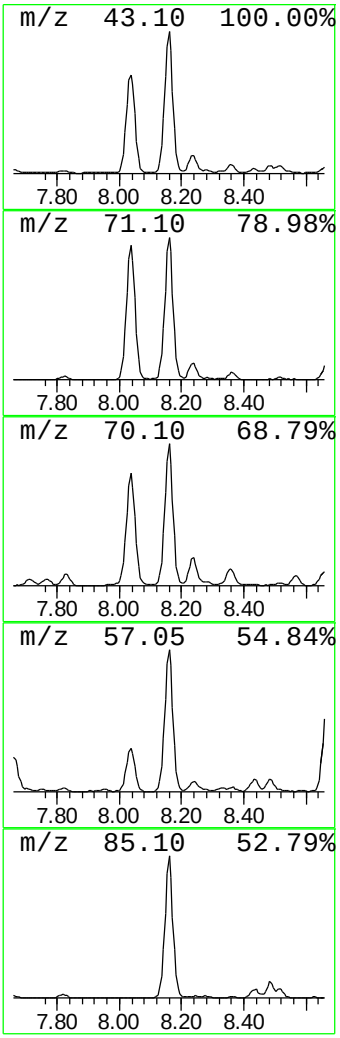
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.16	26.06 ug/l	540959	1,4-Difluorobenzene	6.83

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	74
3	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	72
4	Hexane, 3-methyl-	100	C7H16	000589-34-4	59
5	Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	59





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX102220\
Data File : VX019033.D
Acq On : 22 Oct 2020 17:30
Operator : JC/SP
Sample : L4417-10
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-14

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102120W.M
Quant Title : SW846 8260

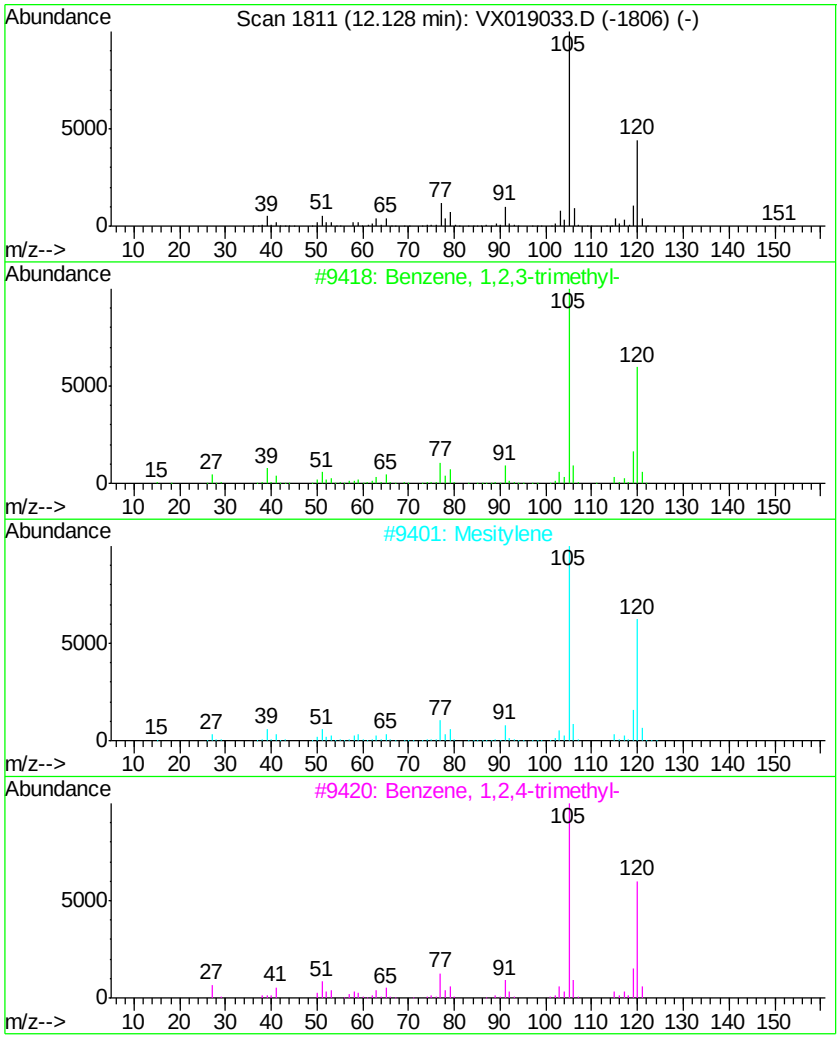
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

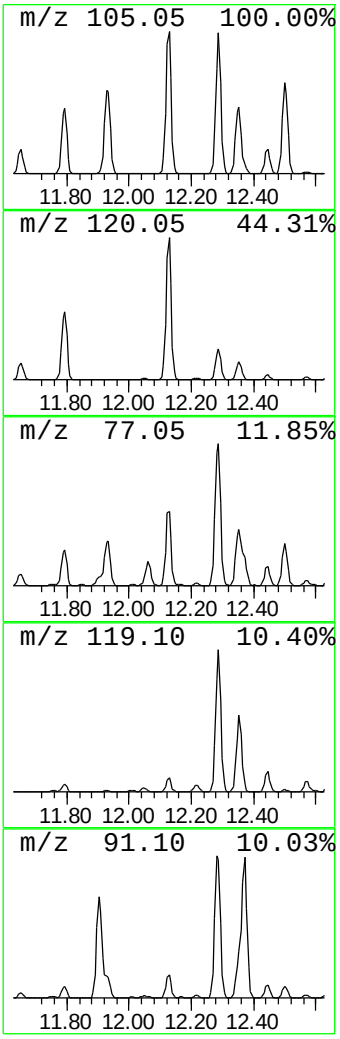
****

Peak Number 2 Benzene, 1,2,3-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.13	25.94 ug/l	738814	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2	Mesitylene	120	C9H12	000108-67-8	94
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX102220\
Data File : VX019033.D
Acq On : 22 Oct 2020 17:30
Operator : JC/SP
Sample : L4417-10
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-14

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102120W.M
Quant Title : SW846 8260

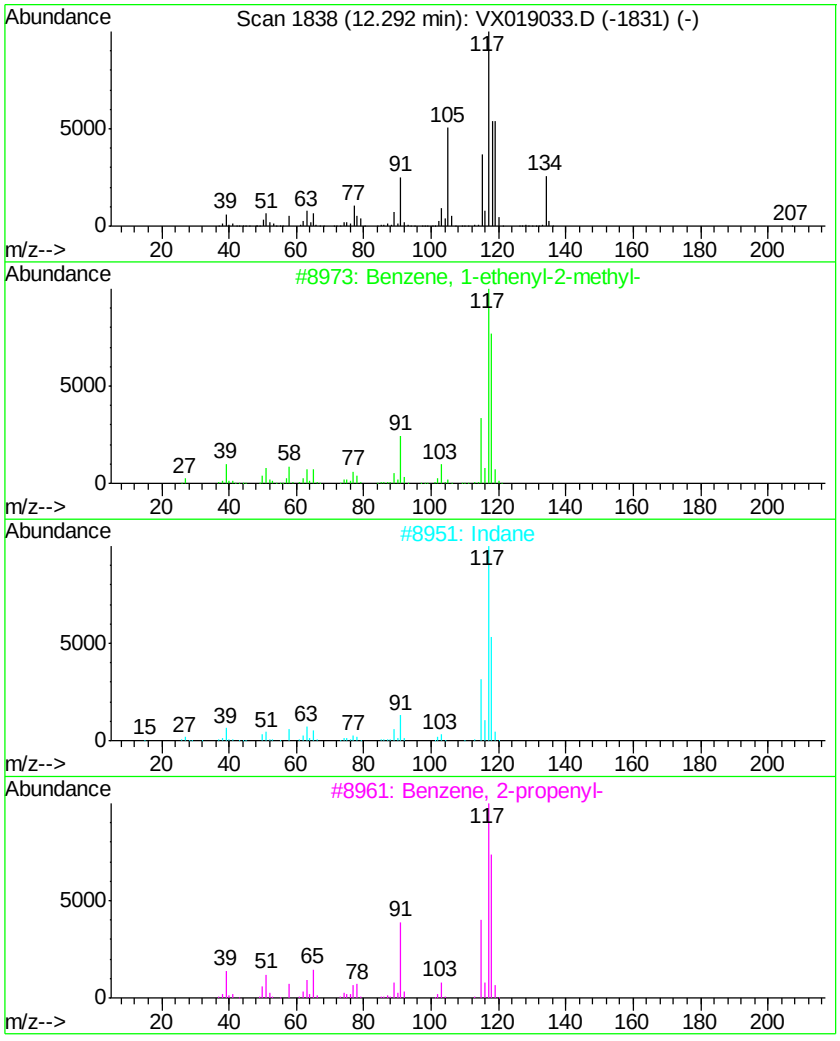
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

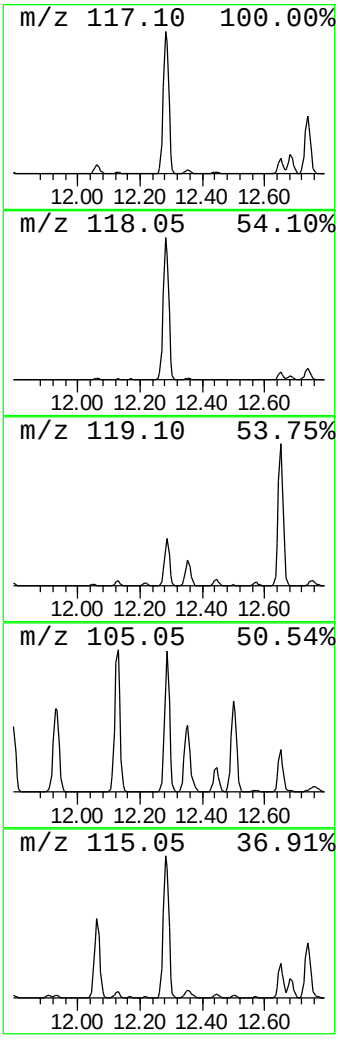
****

Peak Number 3 Benzene, 1-ethenyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	111.19 ug/l	3166890	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	90
2	Indane	118	C9H10	000496-11-7	60
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	60
4	Deltacyclene	118	C9H10	007785-10-6	60
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	60





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX102220\
Data File : VX019033.D
Acq On : 22 Oct 2020 17:30
Operator : JC/SP
Sample : L4417-10
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-14

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102120W.M
Quant Title : SW846 8260

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

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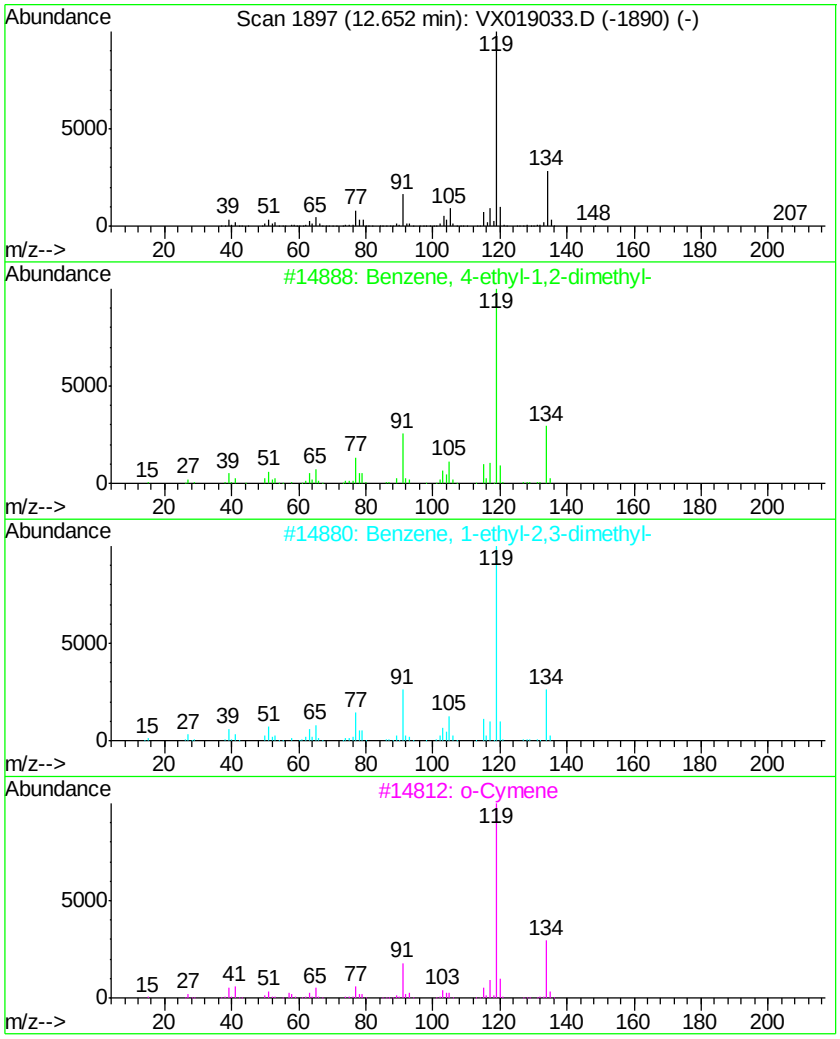
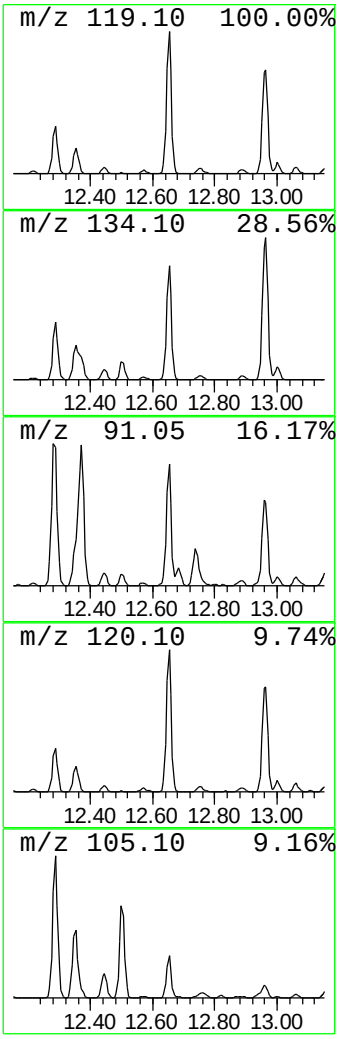
Peak Number 4 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	74.30 ug/l	2116270	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
3	o-Cymene	134	C10H14	000527-84-4	95
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX102220\
Data File : VX019033.D
Acq On : 22 Oct 2020 17:30
Operator : JC/SP
Sample : L4417-10
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-14

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102120W.M
Quant Title : SW846 8260

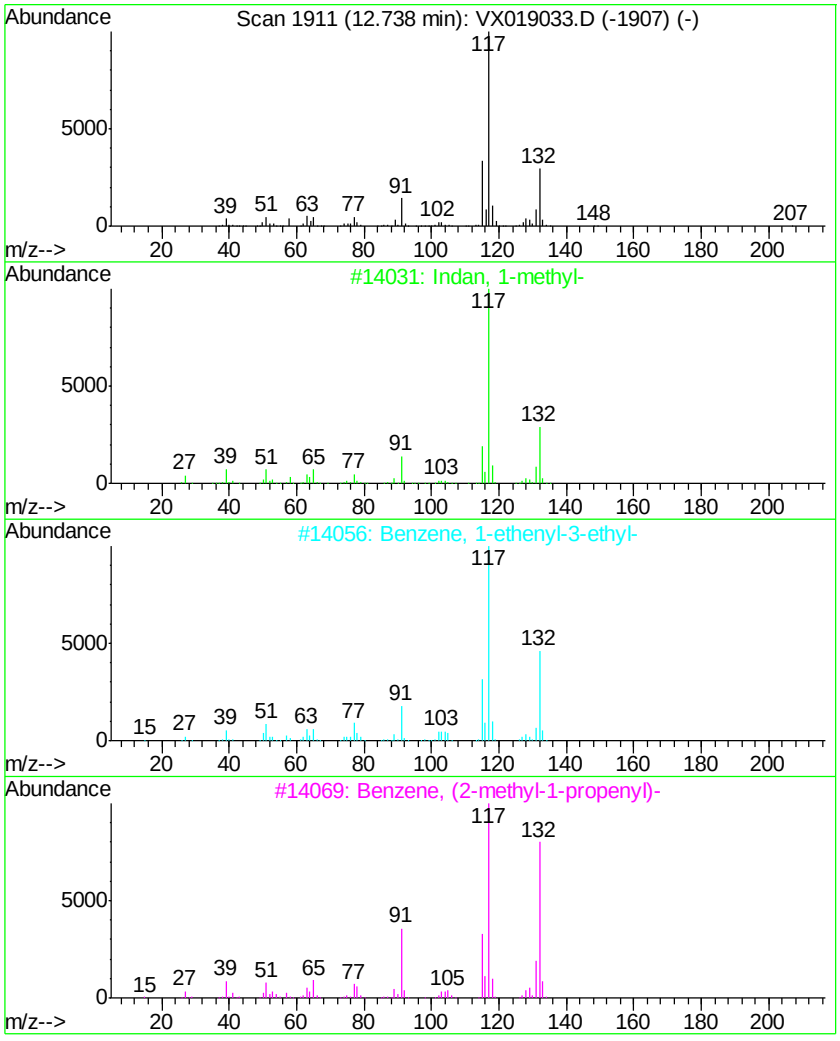
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

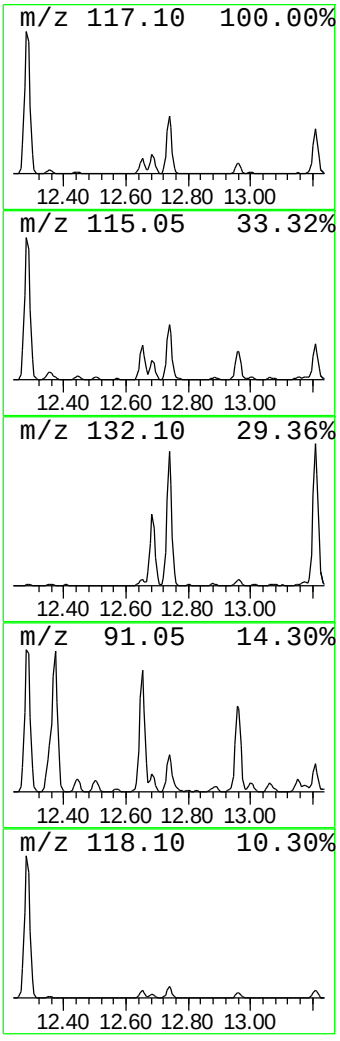
****

Peak Number 5 Indan, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.74	33.14 ug/l	943790	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	91
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
3	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	90
4	Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	87
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	86





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX102220\
Data File : VX019033.D
Acq On : 22 Oct 2020 17:30
Operator : JC/SP
Sample : L4417-10
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-14

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102120W.M
Quant Title : SW846 8260

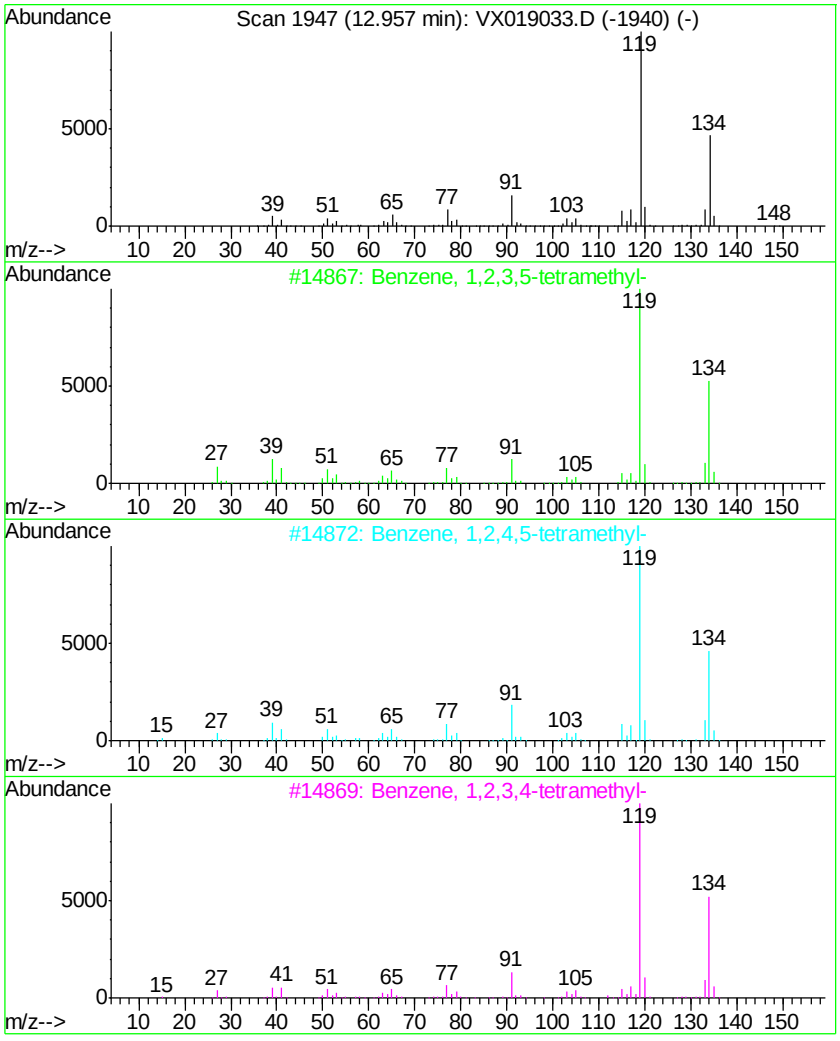
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

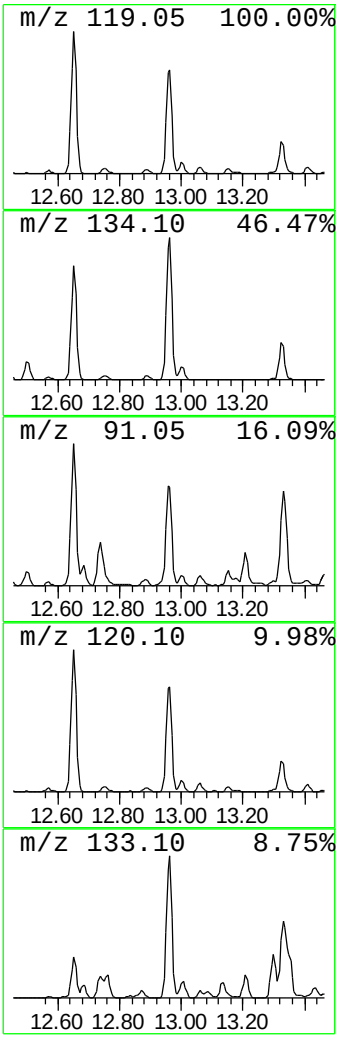
****

Peak Number 6 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.96	64.66 ug/l	1841670	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94





Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX102220\
Data File : VX019033.D
Acq On : 22 Oct 2020 17:30
Operator : JC/SP
Sample : L4417-10
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-14

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102120W.M
Quant Title : SW846 8260

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

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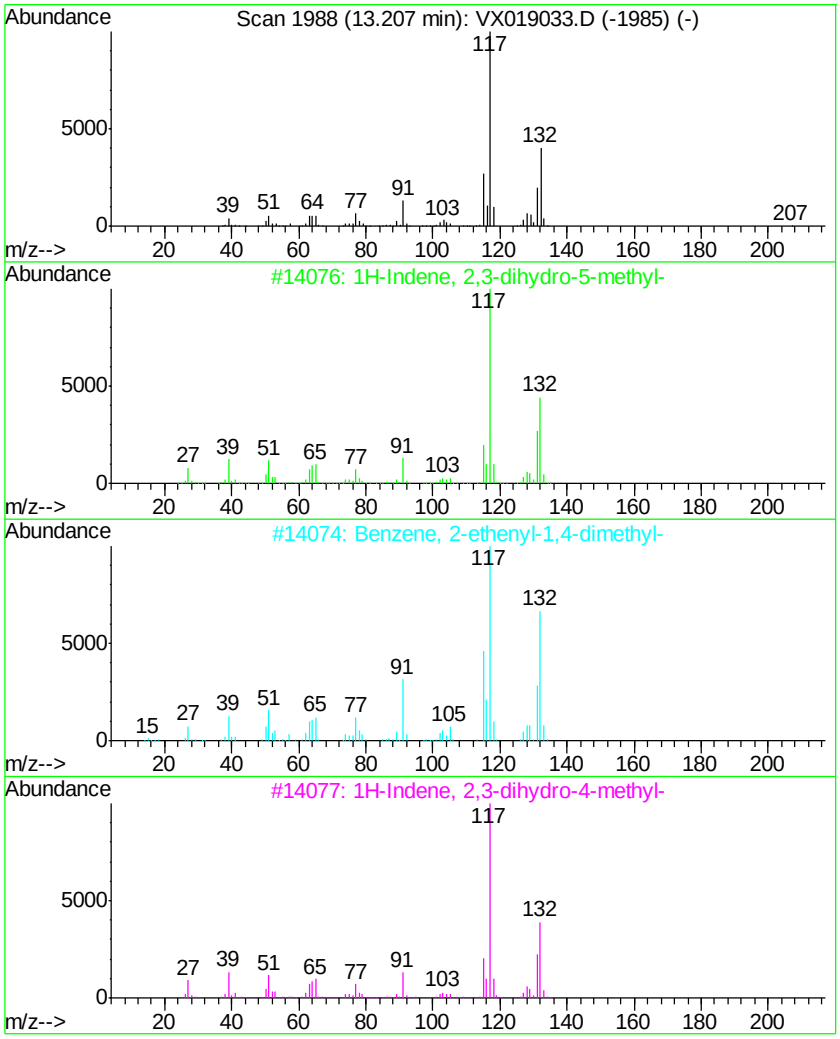
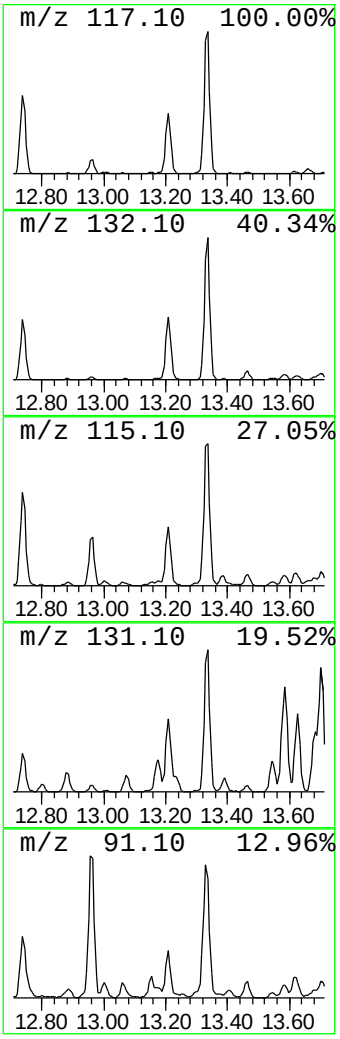
Peak Number 7 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.21	25.28 ug/l	719991	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
2	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	92
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
5	Indan, 1-methyl-	132	C10H12	000767-58-8	87





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX102220\
Data File : VX019033.D
Acq On : 22 Oct 2020 17:30
Operator : JC/SP
Sample : L4417-10
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-14

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102120W.M
Quant Title : SW846 8260

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

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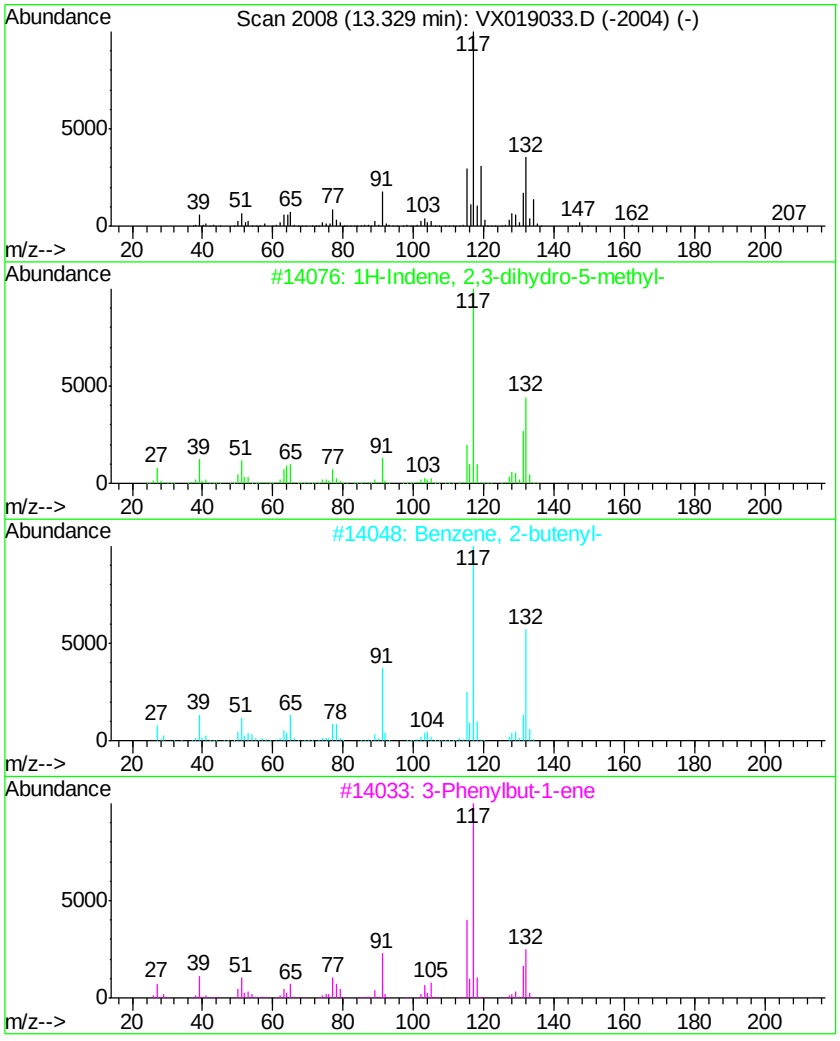
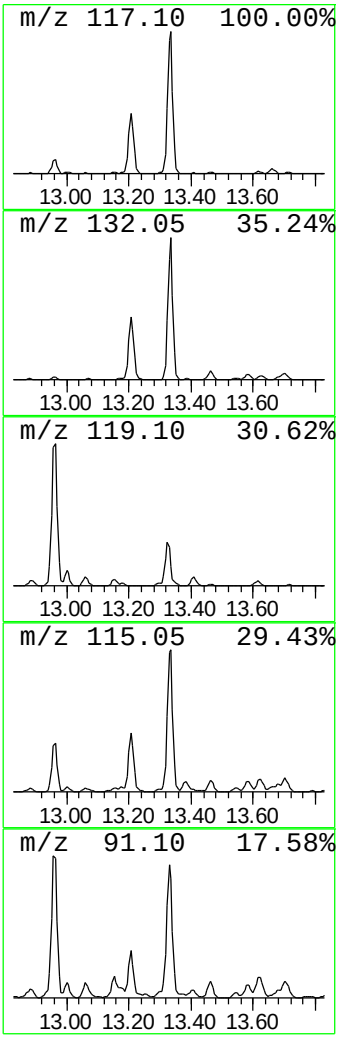
Peak Number 8 Benzene, 2-butenyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.33	74.96 ug/l	2135120	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	89
2	Benzene, 2-butenyl-	132	C10H12	001560-06-1	87
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87
5	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	83





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX102220\
Data File : VX019033.D
Acq On : 22 Oct 2020 17:30
Operator : JC/SP
Sample : L4417-10
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-14

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102120W.M
Quant Title : SW846 8260

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

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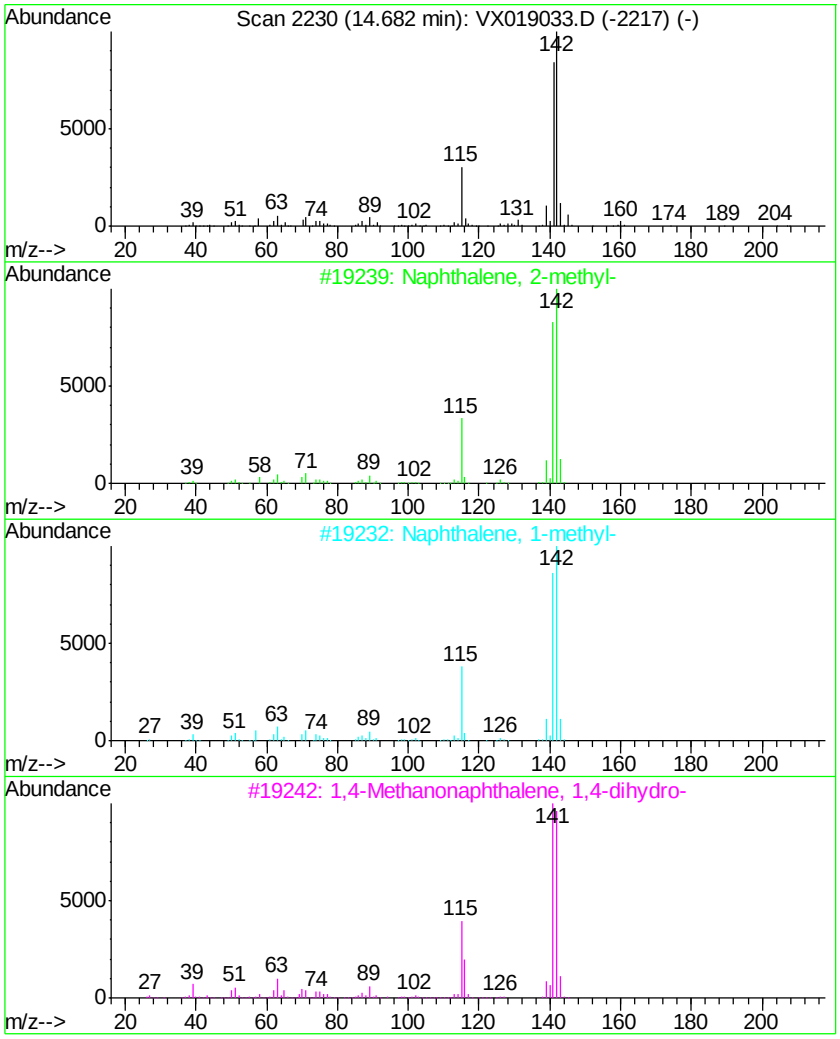
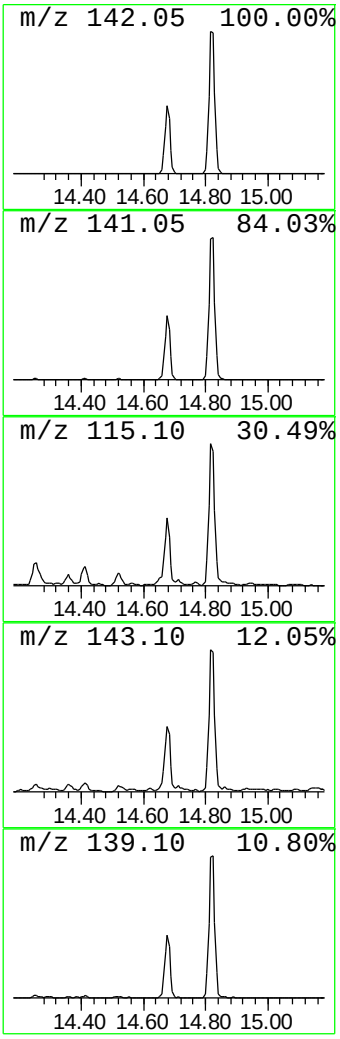
Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.68	14.53 ug/l	413824	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
4	Benzocycloheptatriene	142	C11H10	000264-09-5	91
5	Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	87





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX102220\
Data File : VX019033.D
Acq On : 22 Oct 2020 17:30
Operator : JC/SP
Sample : L4417-10
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-14

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102120W.M
Quant Title : SW846 8260

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

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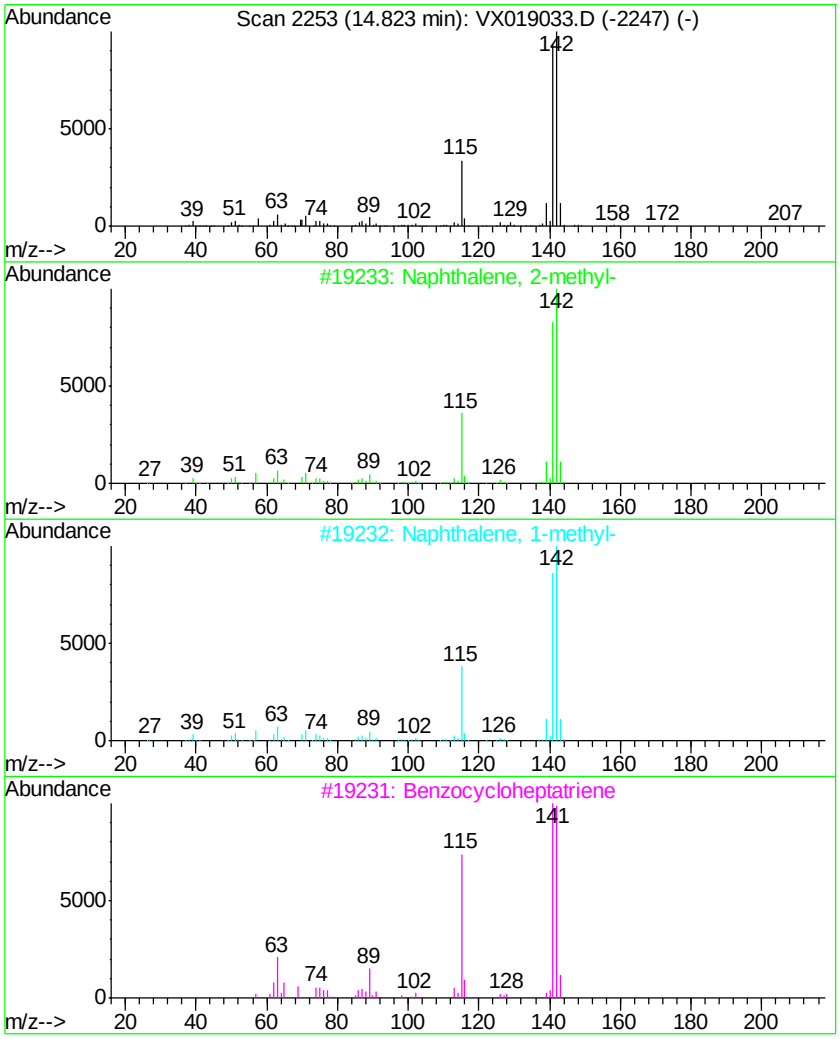
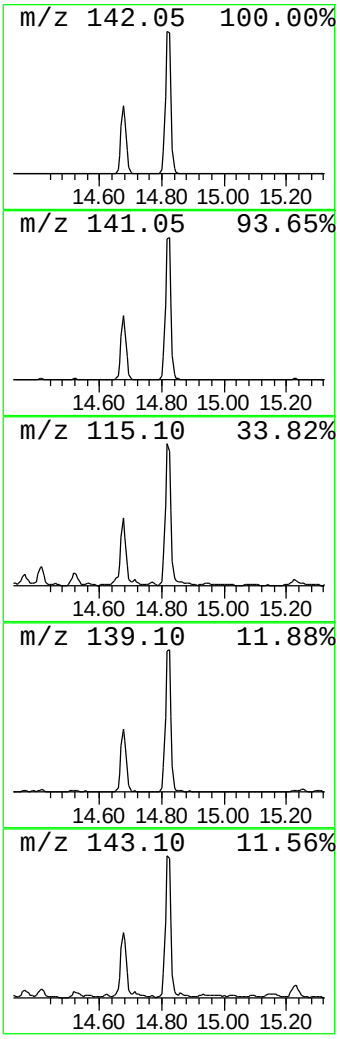
Peak Number 10 Benzocycloheptatriene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	24.91 ug/l	709551	1,4-Dichlorobenzene-d4	12.06

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3	Benzocycloheptatriene	142	C11H10	000264-09-5	91
4	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX102220\
Data File : VX019033.D
Acq On : 22 Oct 2020 17:30
Operator : JC/SP
Sample : L4417-10
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument : MSVOA_X
ClientSampleId : MW-14

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X102120W.M
Quant Title : SW846 8260

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

					Internal Standard		
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Pentane, 2,3,3-tr...	8.16	26.1	ug/l	540959	2	6.83	1037860	50.0
Benzene, 1,2,3-tr...	12.13	25.9	ug/l	738814	4	12.06	1424160	50.0
Benzene, 1-etheny...	12.29	111.2	ug/l	3166890	4	12.06	1424160	50.0
Benzene, 4-ethyl-...	12.65	74.3	ug/l	2116270	4	12.06	1424160	50.0
Indan, 1-methyl-	12.74	33.1	ug/l	943790	4	12.06	1424160	50.0
Benzene, 1,2,3,5-...	12.96	64.7	ug/l	1841670	4	12.06	1424160	50.0
1H-Indene, 2,3-di...	13.21	25.3	ug/l	719991	4	12.06	1424160	50.0
Benzene, 2-butenyl-	13.33	75.0	ug/l	2135120	4	12.06	1424160	50.0
Naphthalene, 1-me...	14.68	14.5	ug/l	413824	4	12.06	1424160	50.0
Benzocycloheptatr...	14.82	24.9	ug/l	709551	4	12.06	1424160	50.0