

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX052219\
 Data File : VX009840.D
 Acq On : 23 May 2019 06:21
 Operator : JC/SP
 Sample : K2976-21 50X
 Misc : 6.62g/5mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 EP1-SBR-4BR(14-17)

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % 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:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.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	5.501	701	713	725	rBV2	105640	302026	15.28%	0.710%
2	5.659	727	739	756	rVV	187769	536791	27.16%	1.263%
3	6.062	794	805	819	rBV	133604	351681	17.79%	0.827%
4	6.860	926	936	952	rBV	310743	732861	37.08%	1.724%
5	8.714	1234	1240	1254	rVB	659377	1033600	52.30%	2.431%
6	9.043	1287	1294	1302	rVB	38423	65126	3.30%	0.153%
7	9.165	1305	1314	1319	rBV	17966	30100	1.52%	0.071%
8	9.348	1338	1344	1351	rVB2	21169	35575	1.80%	0.084%
9	9.445	1351	1360	1369	rBV	96667	159369	8.06%	0.375%
10	9.561	1373	1379	1385	rBV2	69094	146955	7.44%	0.346%
11	9.677	1392	1398	1402	rBV	218453	342488	17.33%	0.806%
12	9.719	1402	1405	1408	rVV2	41824	71450	3.62%	0.168%
13	9.750	1408	1410	1414	rVV2	24518	31523	1.59%	0.074%
14	9.817	1414	1421	1425	rVV	46904	78172	3.96%	0.184%
15	9.884	1425	1432	1437	rVV3	149669	371556	18.80%	0.874%
16	9.951	1441	1443	1448	rVB	88617	112737	5.70%	0.265%
17	10.061	1455	1461	1465	rBV	150863	227163	11.49%	0.534%
18	10.110	1465	1469	1474	rVB	657964	868022	43.92%	2.042%
19	10.177	1474	1480	1485	rVB5	38372	90658	4.59%	0.213%
20	10.238	1485	1490	1496	rBV2	87342	148419	7.51%	0.349%
21	10.335	1496	1506	1508	rBV2	175563	341812	17.29%	0.804%
22	10.372	1508	1512	1516	rVV	268647	370503	18.75%	0.871%
23	10.408	1516	1518	1530	rVB	193118	328415	16.62%	0.772%
24	10.512	1530	1535	1541	rBV	133610	200708	10.16%	0.472%
25	10.585	1543	1547	1549	rBV	43617	57138	2.89%	0.134%
26	10.640	1549	1556	1564	rBV3	506598	894704	45.27%	2.105%
27	10.725	1564	1570	1573	rBV3	262870	411844	20.84%	0.969%
28	10.762	1574	1576	1580	rVB2	134663	166893	8.44%	0.393%
29	10.835	1580	1588	1592	rBV	1343649	1976390	100.00%	4.649%
30	10.884	1592	1596	1597	rBV3	131762	176365	8.92%	0.415%
31	10.908	1597	1600	1603	rVV3	368702	515822	26.10%	1.213%
32	10.945	1603	1606	1611	rVB	601242	831471	42.07%	1.956%
33	11.000	1611	1615	1620	rVB2	176889	283612	14.35%	0.667%
34	11.055	1620	1624	1626	rBV2	263326	384849	19.47%	0.905%

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Integrator: RTE
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35	11.085	1626	1629	1633	rVB2	366705	457527	23.15%	1.076%
36	11.134	1633	1637	1642	rVB	1170305	1508375	76.32%	3.548%
37	11.183	1642	1645	1648	rBV	168368	192871	9.76%	0.454%
38	11.219	1648	1651	1653	rBV	123202	150463	7.61%	0.354%
39	11.274	1653	1660	1669	rVB3	920090	1467344	74.24%	3.451%
40	11.390	1675	1679	1684	rBV2	299973	579367	29.31%	1.363%
41	11.445	1684	1688	1691	rBV3	280188	365289	18.48%	0.859%
42	11.481	1691	1694	1698	rBV	724344	887220	44.89%	2.087%
43	11.536	1698	1703	1707	rVB2	456857	770134	38.97%	1.812%
44	11.579	1707	1710	1715	rVB5	136183	178453	9.03%	0.420%
45	11.628	1715	1718	1720	rBV	128167	154692	7.83%	0.364%
46	11.683	1720	1727	1732	rVB4	421235	822682	41.63%	1.935%
47	11.737	1732	1736	1738	rBV2	350384	466902	23.62%	1.098%
48	11.768	1738	1741	1744	rVV	1574070	1643076	83.14%	3.865%
49	11.804	1744	1747	1757	rVB5	557633	1468646	74.31%	3.455%
50	11.890	1757	1761	1763	rBV	301832	429864	21.75%	1.011%
51	11.945	1763	1770	1775	rVB5	678718	1160912	58.74%	2.731%
52	12.000	1775	1779	1781	rBV	640242	852887	43.15%	2.006%
53	12.024	1781	1783	1787	rVB3	696011	740633	37.47%	1.742%
54	12.079	1788	1792	1797	rVB2	782190	1239426	62.71%	2.915%
55	12.128	1797	1800	1803	rVB4	179300	197251	9.98%	0.464%
56	12.164	1803	1806	1808	rBV3	67486	78984	4.00%	0.186%
57	12.213	1812	1814	1823	rVB4	248078	529457	26.79%	1.245%
58	12.323	1823	1832	1835	rBV2	1085425	1932432	97.78%	4.545%
59	12.365	1835	1839	1846	rVB2	1001401	1483201	75.05%	3.489%
60	12.457	1846	1854	1855	rBV6	294888	444667	22.50%	1.046%
61	12.524	1862	1865	1870	rVB4	310805	524987	26.56%	1.235%
62	12.585	1870	1875	1876	rVV	594642	662073	33.50%	1.557%
63	12.609	1876	1879	1885	rVV2	859506	1502296	76.01%	3.534%
64	12.664	1885	1888	1892	rVV	757345	1016608	51.44%	2.391%
65	12.737	1896	1900	1902	rVV3	177950	259675	13.14%	0.611%
66	12.774	1902	1906	1911	rVV5	502725	971051	49.13%	2.284%
67	12.816	1911	1913	1914	rVV2	95430	93152	4.71%	0.219%
68	12.847	1915	1918	1920	rVV2	255798	359198	18.17%	0.845%
69	12.871	1920	1922	1924	rVV	231479	294495	14.90%	0.693%
70	12.902	1924	1927	1930	rVV2	243191	324890	16.44%	0.764%
71	12.944	1930	1934	1937	rVV3	274253	480729	24.32%	1.131%

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Integration Parameters: RTEINT.P

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72	12.975	1937	1939	1942	rVV	328373	353639	17.89%	0.832%
73	13.018	1942	1946	1949	rVV	540711	691920	35.01%	1.628%
74	13.079	1953	1956	1962	rVB3	182103	287032	14.52%	0.675%
75	13.146	1965	1967	1972	rVB2	81279	103592	5.24%	0.244%
76	13.207	1972	1977	1982	rVB4	132793	282668	14.30%	0.665%
77	13.262	1983	1986	1989	rVB2	75092	86380	4.37%	0.203%
78	13.304	1989	1993	1996	rBV2	138727	177272	8.97%	0.417%
79	13.347	1996	2000	2006	rVB4	292212	456096	23.08%	1.073%
80	13.402	2006	2009	2010	rBV3	23030	25305	1.28%	0.060%
81	13.426	2010	2013	2018	rVB2	51513	63207	3.20%	0.149%
82	13.481	2018	2022	2030	rVB3	67225	156035	7.89%	0.367%
83	13.560	2030	2035	2038	rBV3	45134	69277	3.51%	0.163%
84	13.597	2038	2041	2043	rVB	22540	21052	1.07%	0.050%
85	13.633	2043	2047	2053	rBV3	79879	142665	7.22%	0.336%
86	13.719	2059	2061	2066	rVB2	52803	73914	3.74%	0.174%
87	13.804	2072	2075	2079	rVB2	20389	25142	1.27%	0.059%
88	13.853	2079	2083	2086	rVV2	19413	25311	1.28%	0.060%
89	13.908	2086	2092	2097	rVB6	19514	38595	1.95%	0.091%
90	13.981	2097	2104	2109	rVB5	20082	41816	2.12%	0.098%
91	14.267	2147	2151	2159	rVB4	10901	23898	1.21%	0.056%

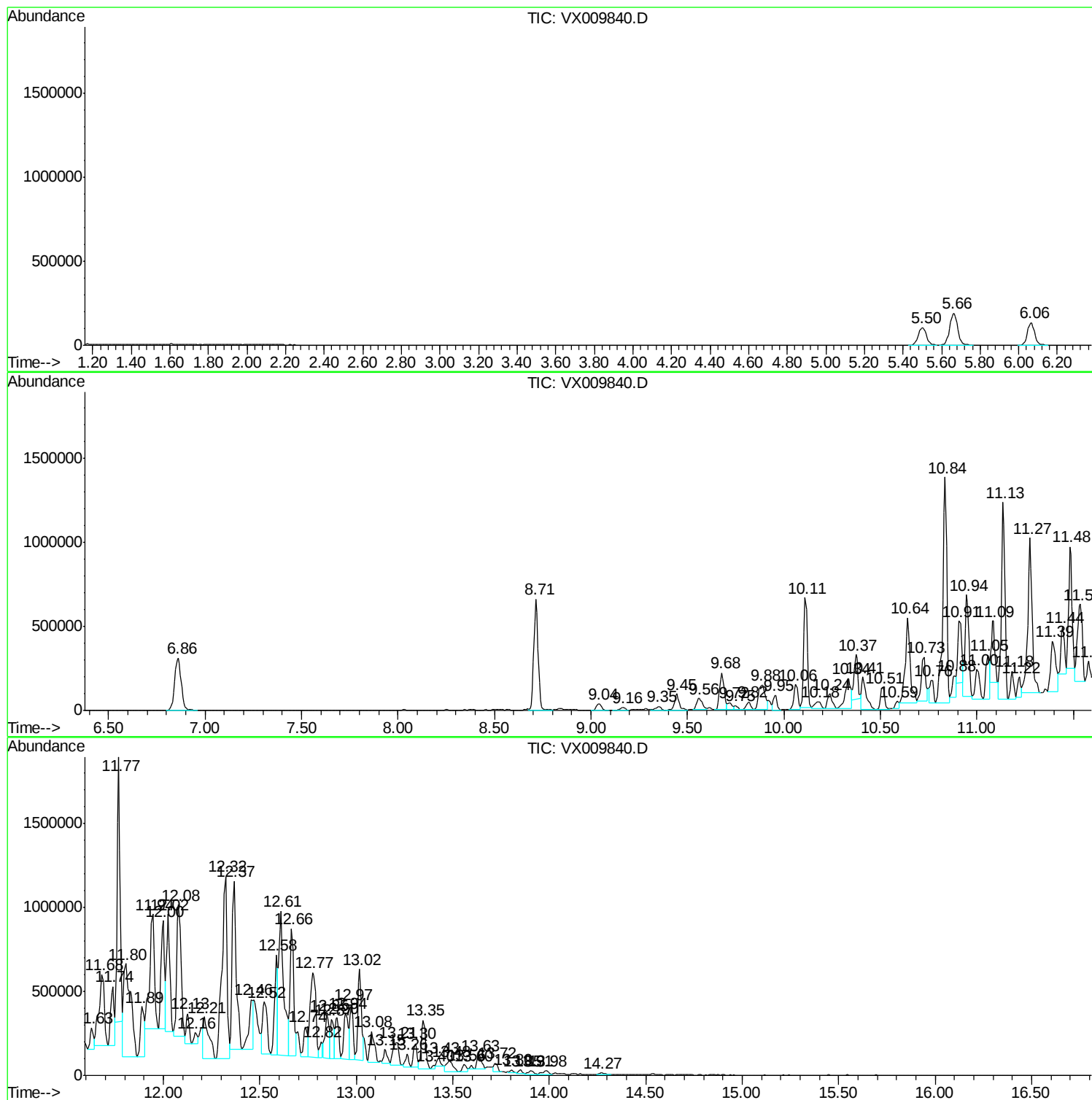
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Instrument :
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ClientSampleId :
EP1-SBR-4BR(14-17)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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



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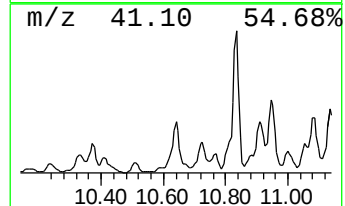
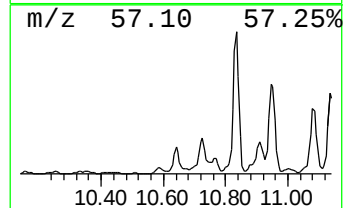
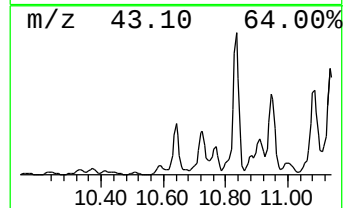
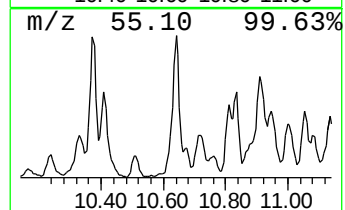
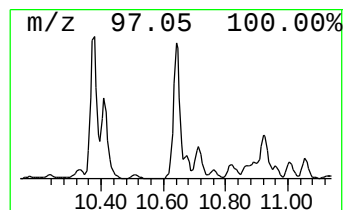
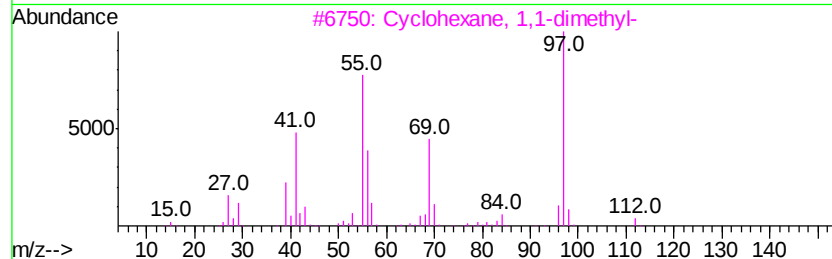
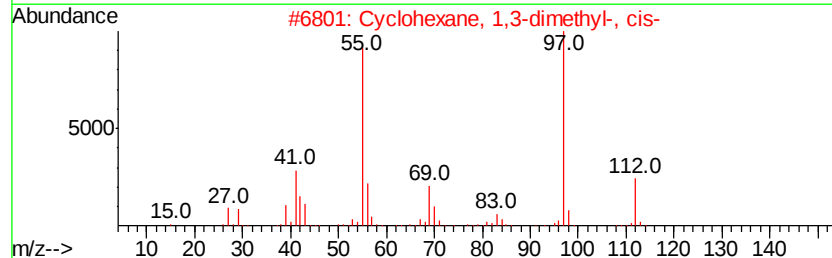
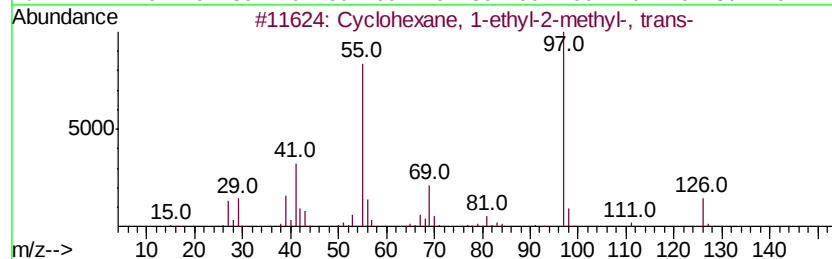
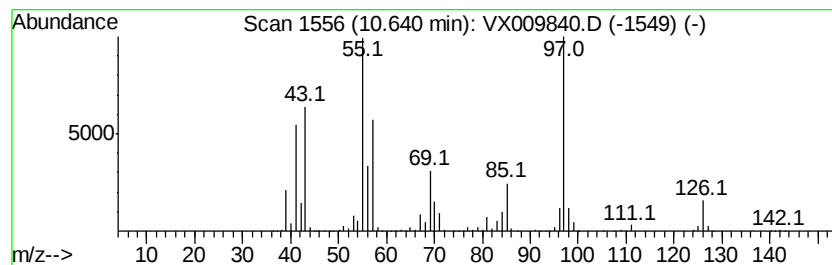
Instrument :
MSVOA_X
ClientSampleId :
EP1-SBR-4BR(14-17)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclohexane, 1-ethyl-2-meth... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.64	51.54 ug/l	894704	Chlorobenzene-d5	10.11		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cvclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	64
2		Cvclohexane, 1,3-dimethyl-, cis-	112	C8H16	000638-04-0	53
3		Cvclohexane, 1,1-dimethyl-	112	C8H16	000590-66-9	53
4		Sulfurous acid, cvclohexylmethyl...	388	C22H44O3S	1000309-22-3	53
5		Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	52



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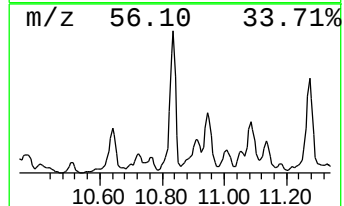
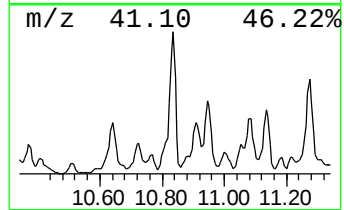
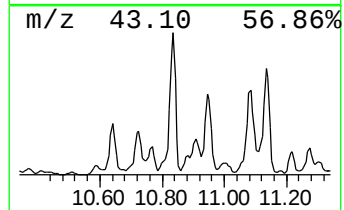
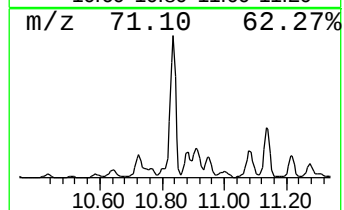
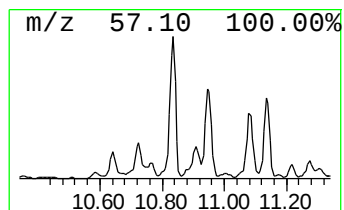
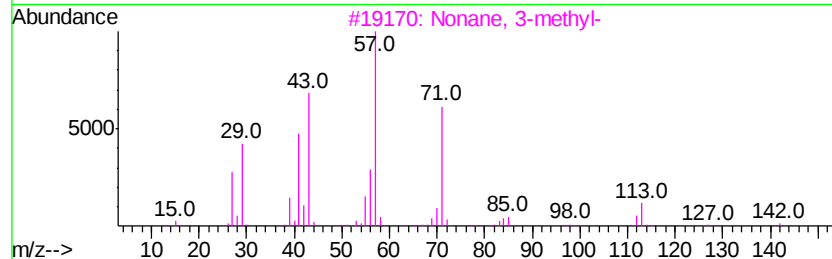
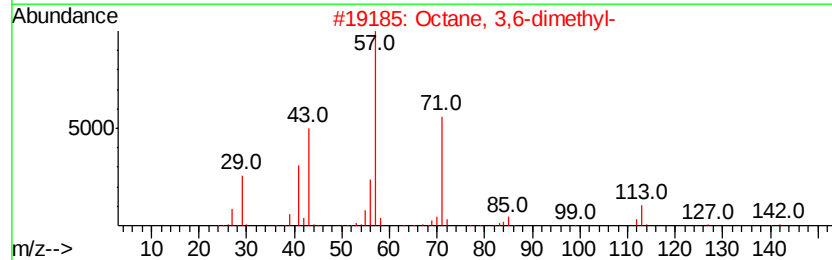
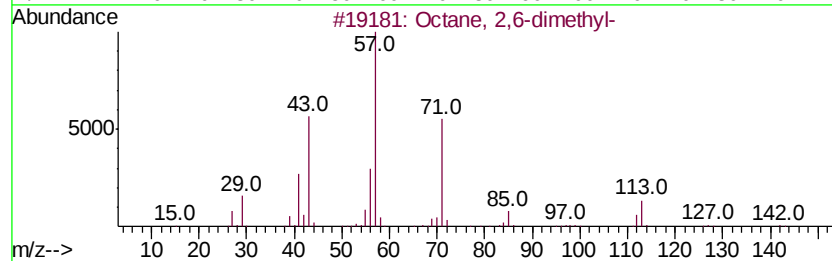
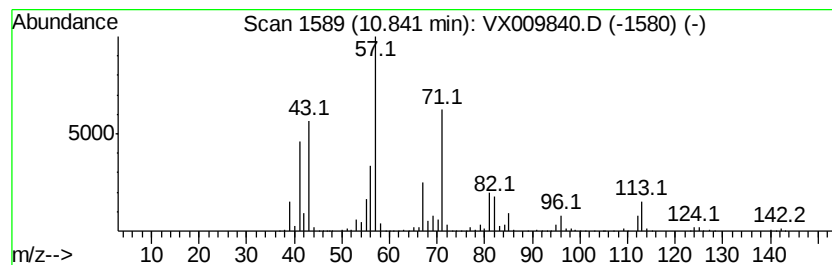
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Octane, 2,6-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.84	113.84 ug/l	1976390	Chlorobenzene-d5	10.11		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	94
2		Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	81
3		Nonane, 3-methyl-	142	C10H22	005911-04-6	76
4		Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	52
5		Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	47



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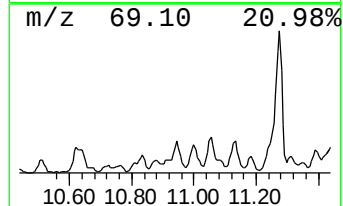
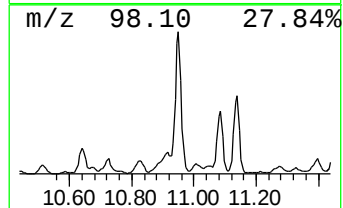
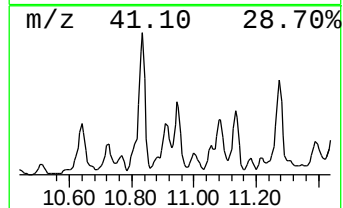
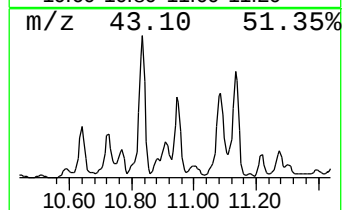
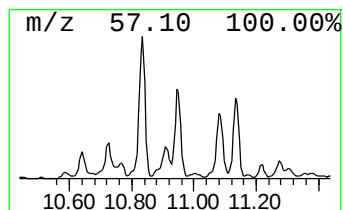
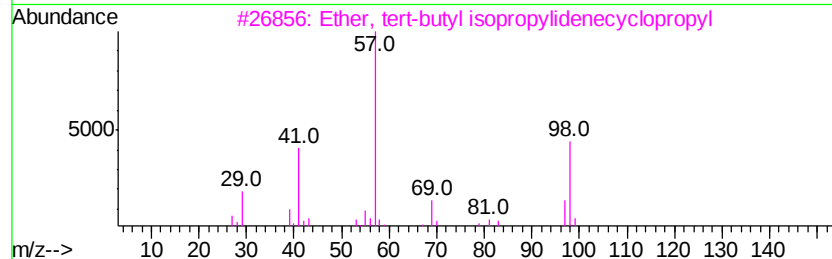
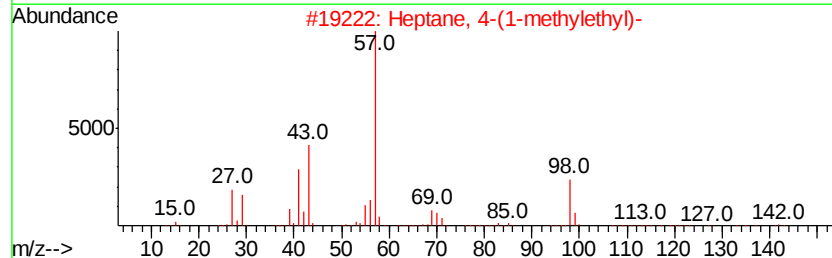
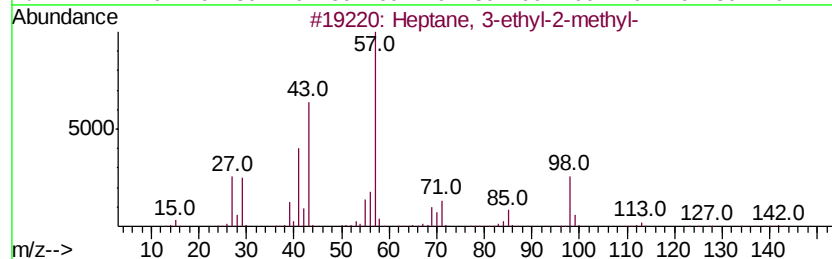
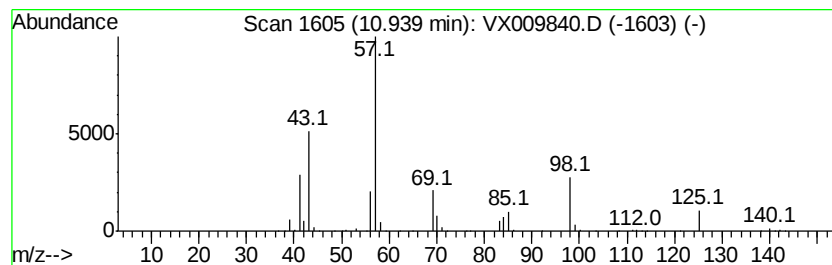
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TIC Integration Parameters: LSCINT.P

Peak Number 3 Heptane, 3-ethyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.94	47.89 ug/l	831471	Chlorobenzene-d5	10.11		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	72	
2	Heptane, 4-(1-methylethyl)-	142	C10H22	052896-87-4	64	
3	Ether, tert-butyl isopropylidene...	154	C10H18O	024524-56-9	50	
4	Hexane, 3-ethyl-2,5-dimethyl-	142	C10H22	052897-04-8	47	
5	Decane, 2,5-dimethyl-	170	C12H26	017312-50-4	43	



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009840.D
Acq On : 23 May 2019 06:21
Operator : JC/SP
Sample : K2976-21 50X
Misc : 6.62g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 52 Sample Multiplier: 1

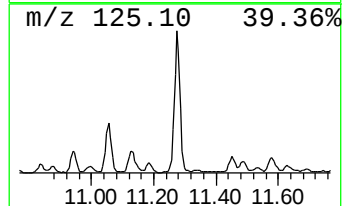
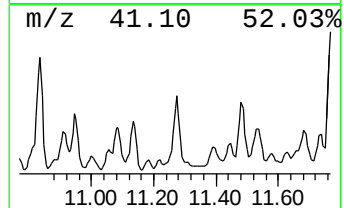
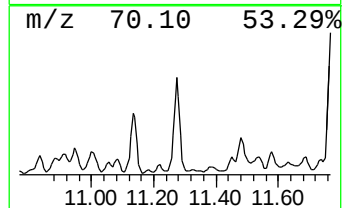
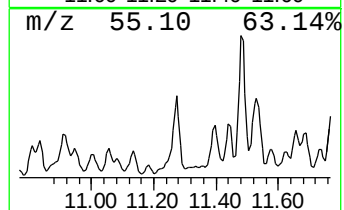
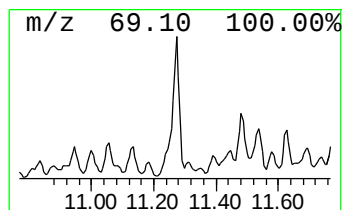
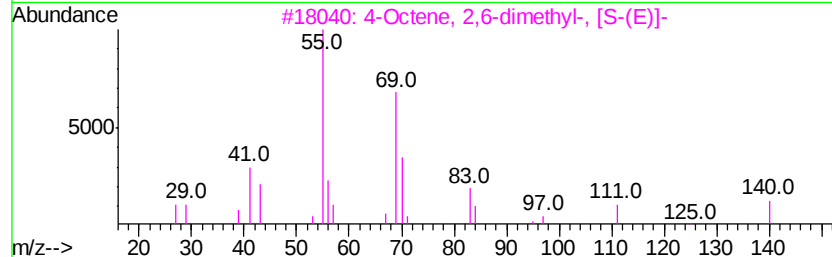
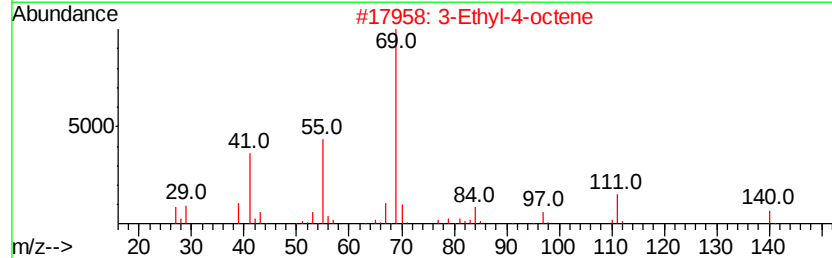
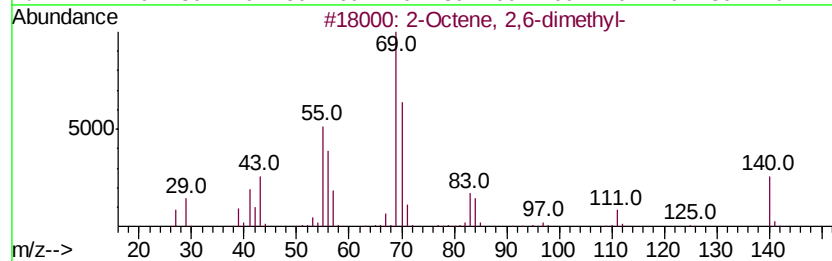
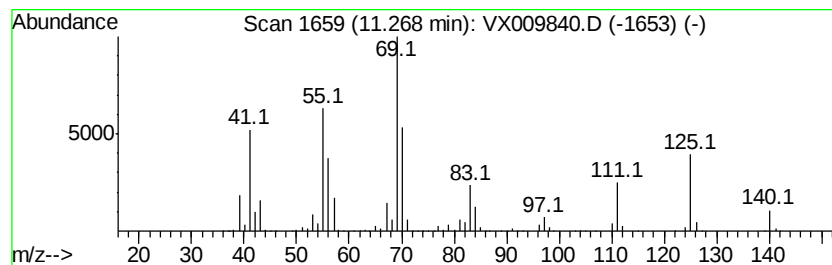
Instrument :
MSVOA_X
ClientSampleId :
EP1-SBR-4BR(14-17)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

Peak Number 4 2-Octene, 2,6-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.27	59.19 ug/l	1467340	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2-Octene, 2,6-dimethyl-	140 C10H20	004057-42-5	64
2	3-Ethyl-4-octene	140 C10H20	019781-32-9	50
3	4-Octene, 2,6-dimethyl-, [S-(E)]-	140 C10H20	062960-76-3	47
4	2-Nonene, 2-methyl-	140 C10H20	002129-95-5	46
5	2-Octene, 2,7-dimethyl-	140 C10H20	002050-81-9	46



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009840.D
Acq On : 23 May 2019 06:21
Operator : JC/SP
Sample : K2976-21 50X
Misc : 6.62g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 52 Sample Multiplier: 1

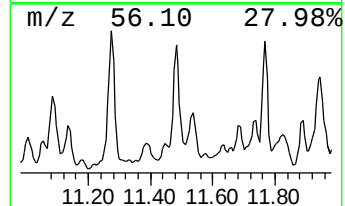
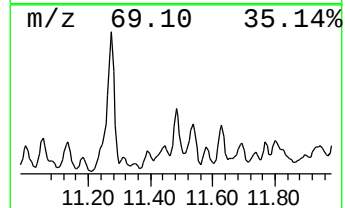
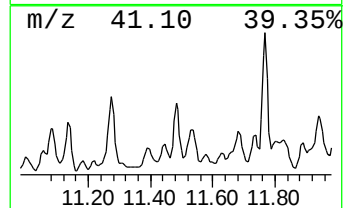
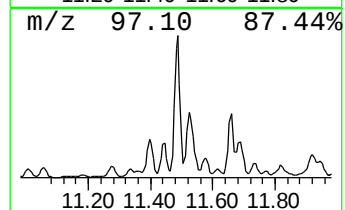
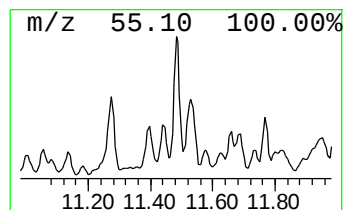
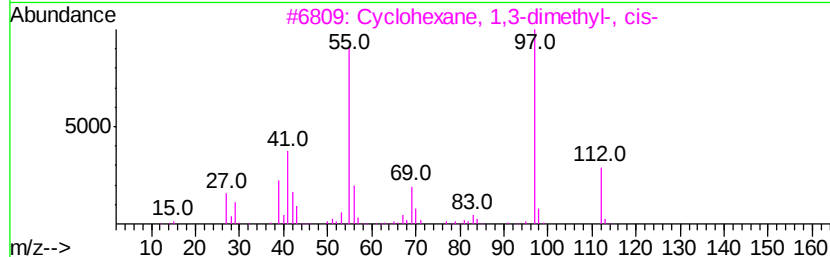
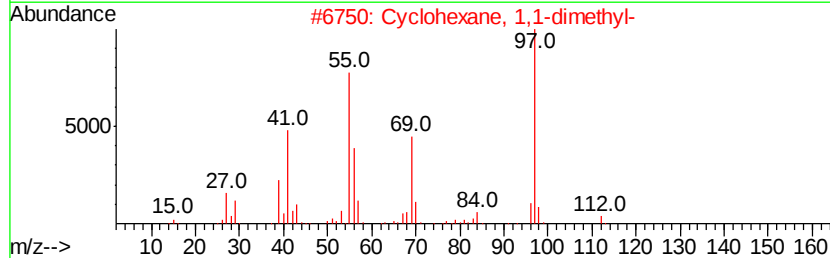
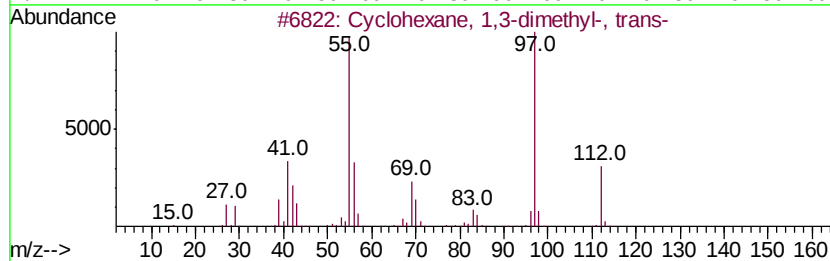
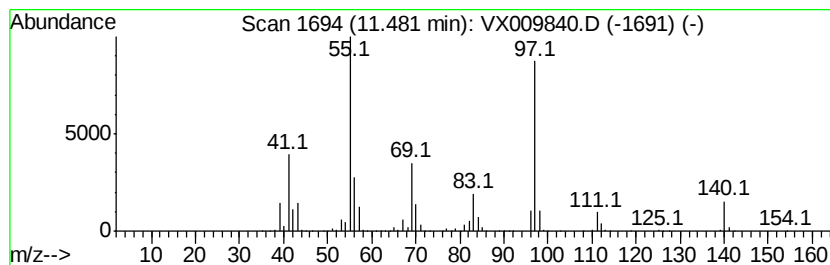
Instrument :
MSVOA_X
ClientSampleId :
EP1-SBR-4BR(14-17)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

Peak Number 5 Cyclohexane, 1,3-dimethyl-,... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.48	35.79 ug/l	887220	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexane, 1,3-dimethyl-, trans-	112 C8H16	002207-03-6 74
2		Cyclohexane, 1,1-dimethyl-	112 C8H16	000590-66-9 70
3		Cyclohexane, 1,3-dimethyl-, cis-	112 C8H16	000638-04-0 68
4		3-Hexene, 3-ethyl-2,5-dimethyl-	140 C10H20	062338-08-3 64
5		Cyclohexane, 1,4-dimethyl-, cis-	112 C8H16	000624-29-3 59



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009840.D
Acq On : 23 May 2019 06:21
Operator : JC/SP
Sample : K2976-21 50X
Misc : 6.62g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 52 Sample Multiplier: 1

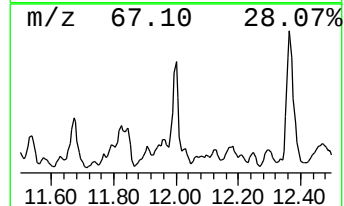
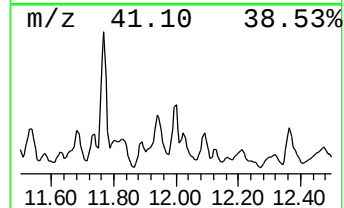
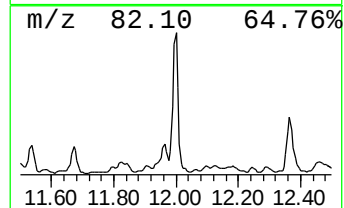
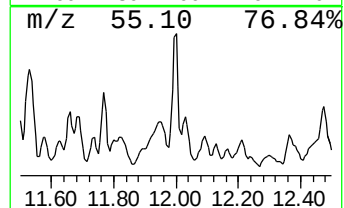
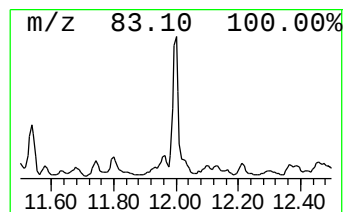
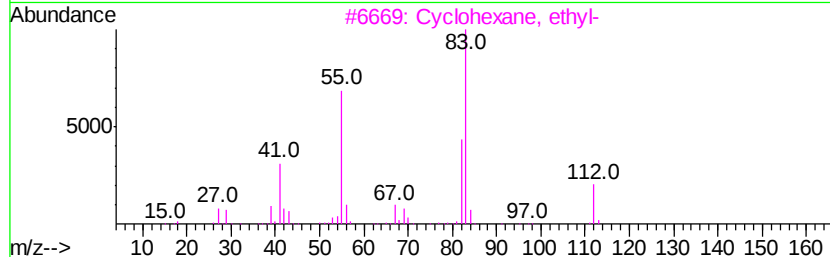
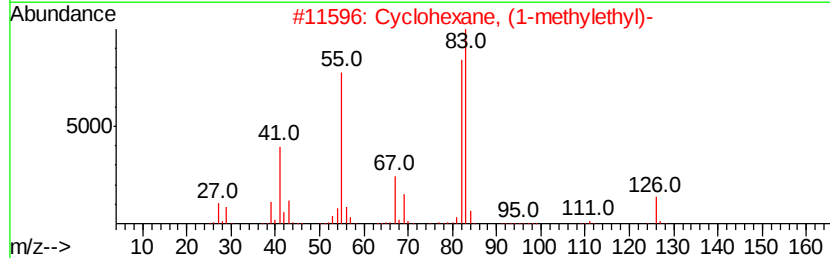
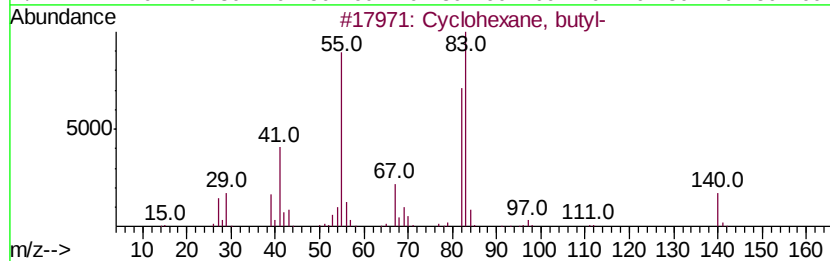
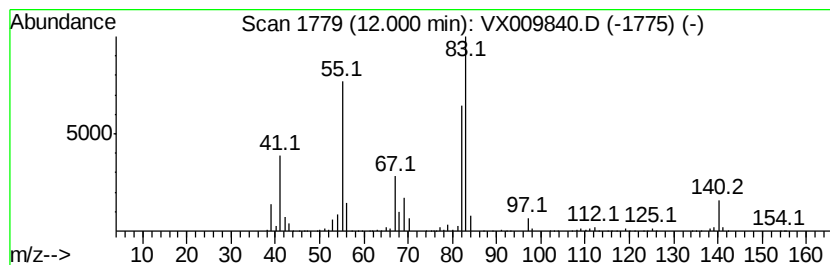
Instrument :
MSVOA_X
ClientSampleId :
EP1-SBR-4BR(14-17)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

Peak Number 6 Cyclohexane, butyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.00	34.41 ug/l	852887	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclohexane, butyl-	140 C10H20	001678-93-9 94
2		Cyclohexane, (1-methylethyl)-	126 C9H18	000696-29-7 72
3		Cyclohexane, ethyl-	112 C8H16	001678-91-7 72
4		Cyclohexane, propyl-	126 C9H18	001678-92-8 72
5		Cyclohexane, pentyl-	154 C11H22	004292-92-6 72



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009840.D
Acq On : 23 May 2019 06:21
Operator : JC/SP
Sample : K2976-21 50X
Misc : 6.62g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 52 Sample Multiplier: 1

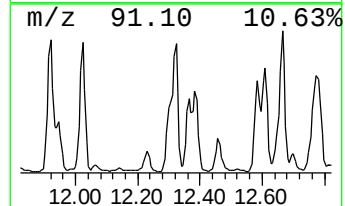
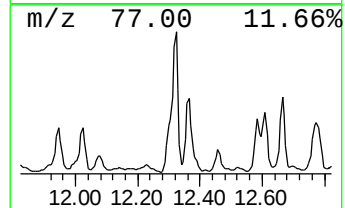
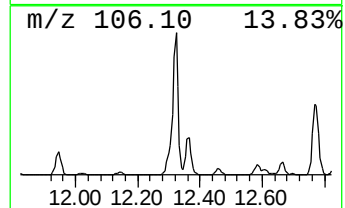
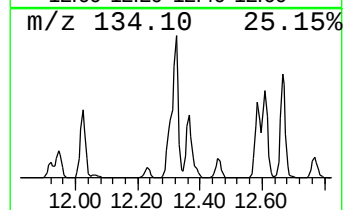
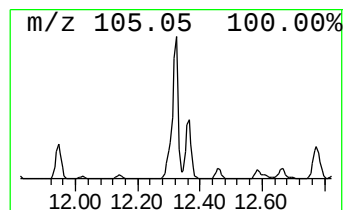
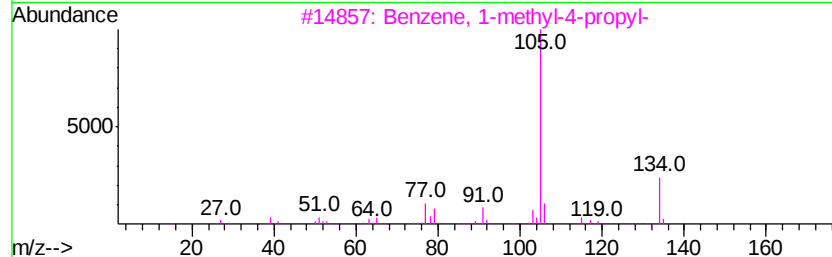
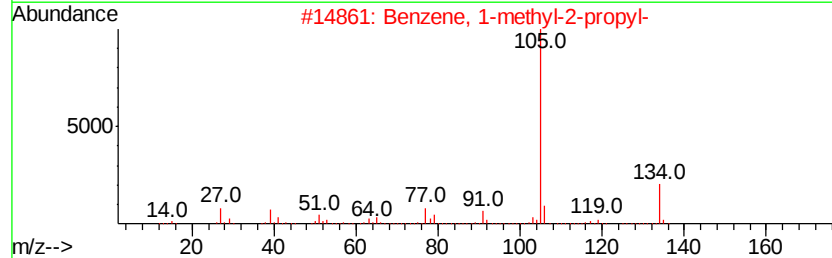
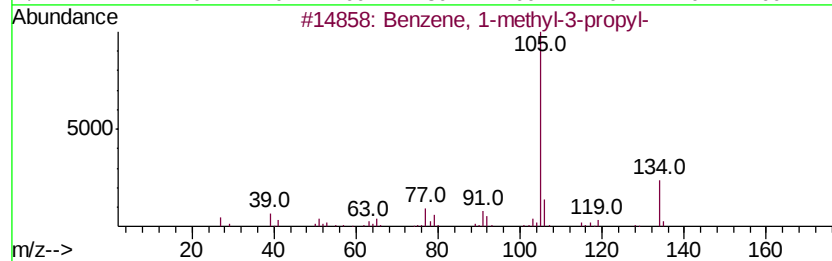
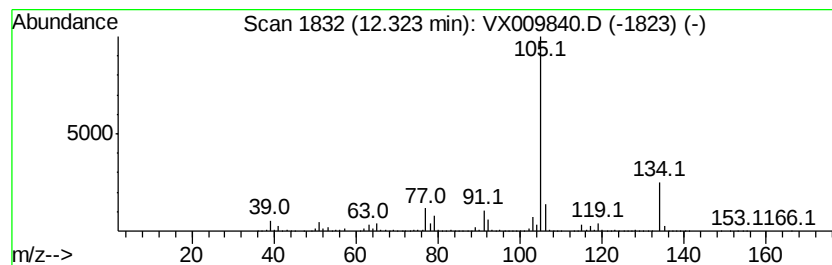
Instrument :
MSVOA_X
ClientSampleId :
EP1-SBR-4BR(14-17)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-methyl-3-propyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.32	77.96 ug/l	1932430	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-methyl-3-propyl-	134 C10H14	001074-43-7 95
2		Benzene, 1-methyl-2-propyl-	134 C10H14	001074-17-5 91
3		Benzene, 1-methyl-4-propyl-	134 C10H14	001074-55-1 91
4		Benzene, (1-methylpropyl)-	134 C10H14	000135-98-8 87
5		2-Tolyloxirane	134 C9H10O	002783-26-8 83



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009840.D
Acq On : 23 May 2019 06:21
Operator : JC/SP
Sample : K2976-21 50X
Misc : 6.62g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 52 Sample Multiplier: 1

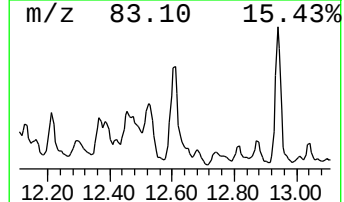
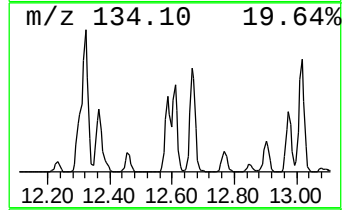
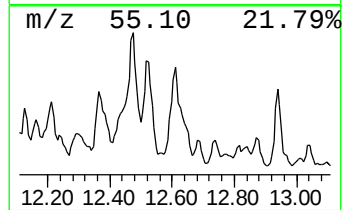
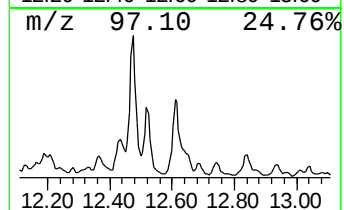
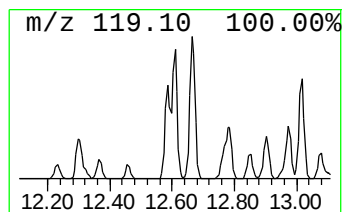
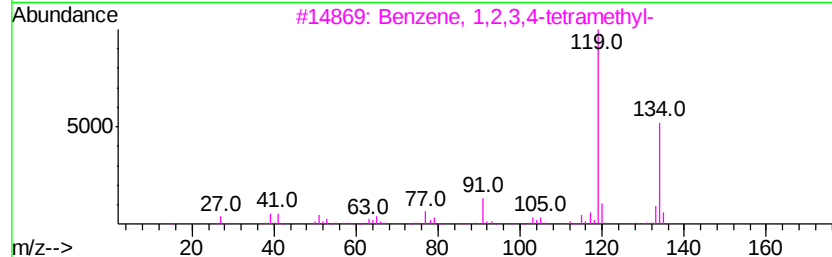
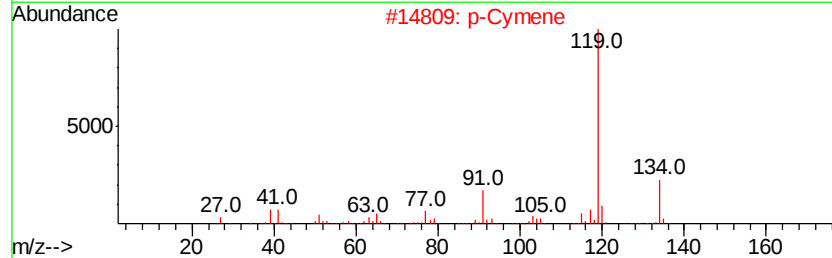
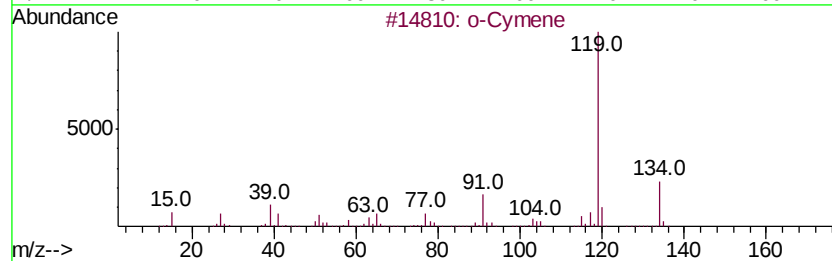
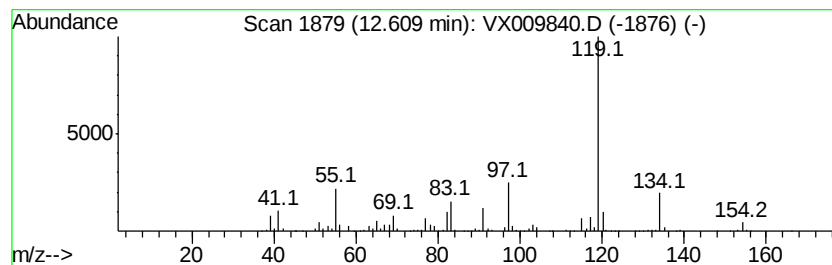
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ClientSampleId :
EP1-SBR-4BR(14-17)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

Peak Number 8 o-Cymene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.61	60.60 ug/l	1502300	1,4-Dichlorobenzene-d4	12.08	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	93
2	p-Cymene	134	C10H14	000099-87-6	92
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	70
4	Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	70
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009840.D
Acq On : 23 May 2019 06:21
Operator : JC/SP
Sample : K2976-21 50X
Misc : 6.62g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 52 Sample Multiplier: 1

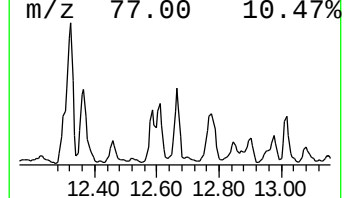
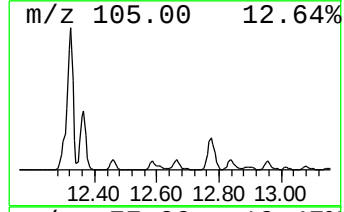
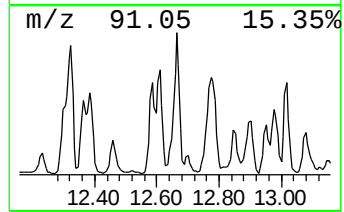
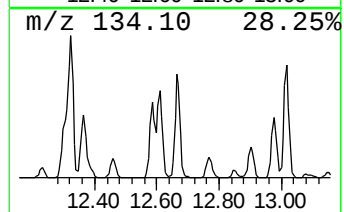
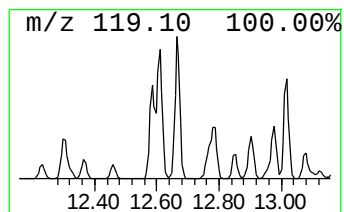
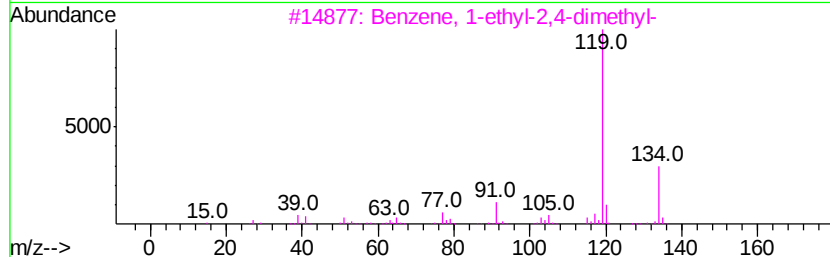
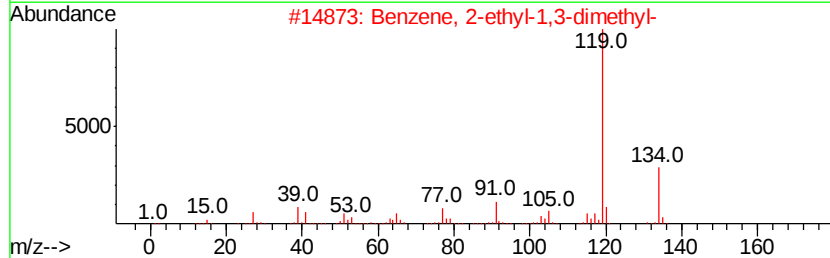
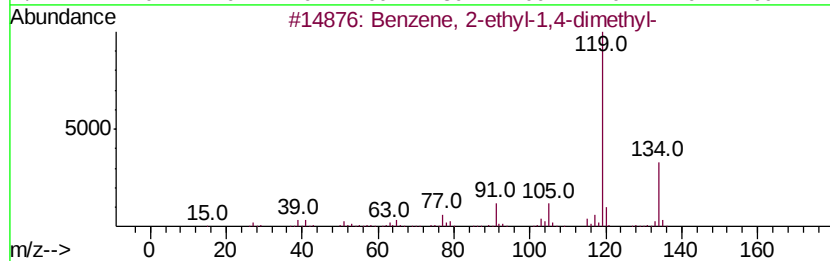
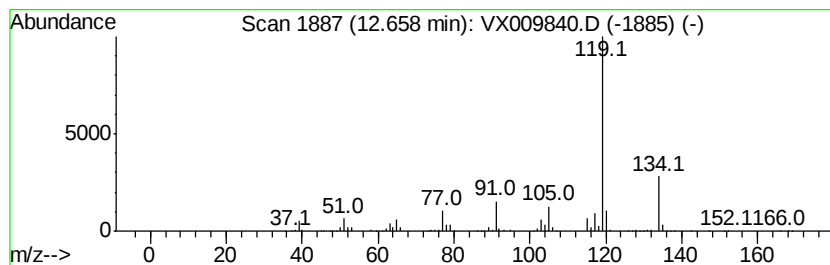
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ClientSampleId :
EP1-SBR-4BR(14-17)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

Peak Number 9 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.66	41.01 ug/l	1016610	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 2-ethyl-1,4-dimethyl-	134 C10H14	001758-88-9 95
2		Benzene, 2-ethyl-1,3-dimethyl-	134 C10H14	002870-04-4 95
3		Benzene, 1-ethyl-2,4-dimethyl-	134 C10H14	000874-41-9 95
4		Benzene, 1-ethyl-2,3-dimethyl-	134 C10H14	000933-98-2 95
5		Benzene, 4-ethyl-1,2-dimethyl-	134 C10H14	000934-80-5 95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX052219\
Data File : VX009840.D
Acq On : 23 May 2019 06:21
Operator : JC/SP
Sample : K2976-21 50X
Misc : 6.62g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 52 Sample Multiplier: 1

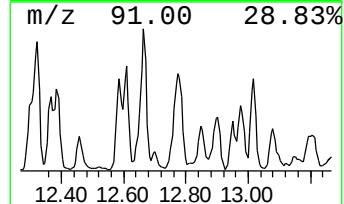
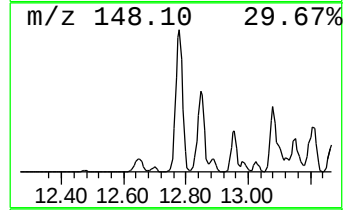
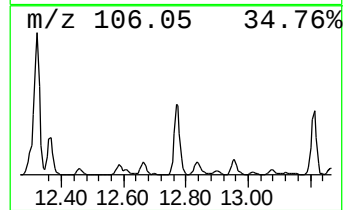
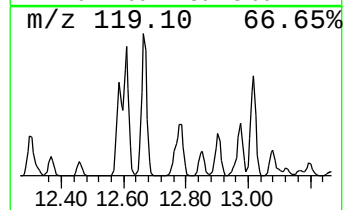
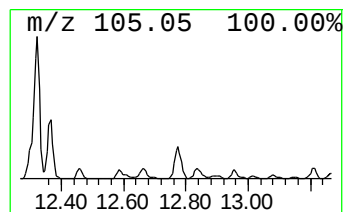
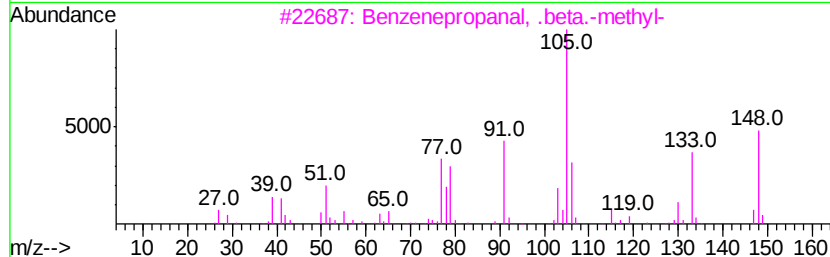
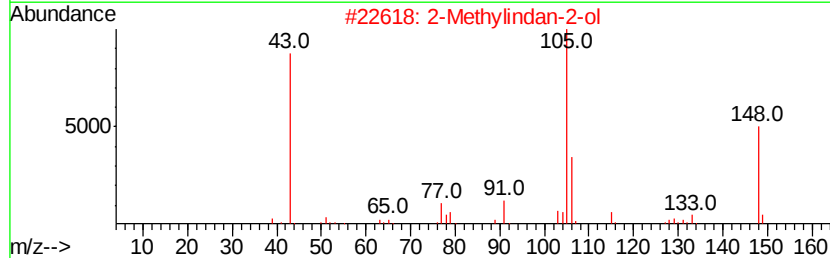
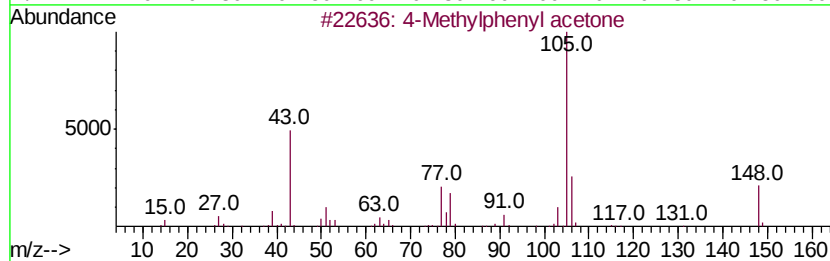
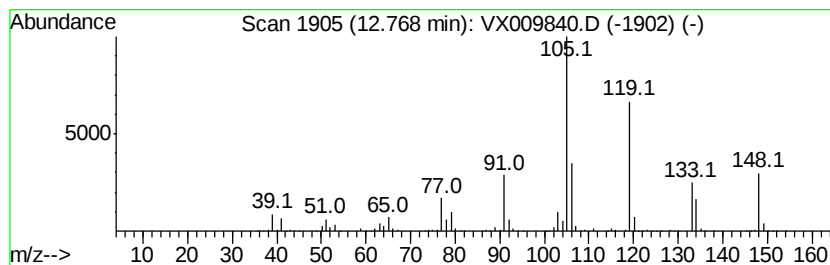
Instrument :
MSVOA_X
ClientSampleId :
EP1-SBR-4BR(14-17)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

Peak Number 10 4-Methylphenyl acetone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.77	39.17 ug/l	971051	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		4-Methylphenyl acetone	148 C10H12O	002096-86-8 55
2		2-Methylindan-2-ol	148 C10H12O	033223-84-6 53
3		Benzenepropanal, .beta.-methyl-	148 C10H12O	016251-77-7 52
4		Benzene, 1-methyl-4-(2-methylpro...	148 C11H16	005161-04-6 49
5		2-Ethyl-2-methyl-1,3-dithiolane	148 C6H12S2	006008-81-7 46



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_X\DATA\VX052219\
Data File : VX009840.D
Acq On : 23 May 2019 06:21
Operator : JC/SP
Sample : K2976-21 50X
Misc : 6.62g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EP1-SBR-4BR(14-17)

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X042919W.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Cyclohexane, 1-et...	10.64	51.5	ug/l	894704	3	10.11	868022	50.0
Octane, 2,6-dimet...	10.84	113.8	ug/l	1976390	3	10.11	868022	50.0
Heptane, 3-ethyl...	10.94	47.9	ug/l	831471	3	10.11	868022	50.0
2-Octene, 2,6-dim...	11.27	59.2	ug/l	1467340	4	12.08	1239430	50.0
Cyclohexane, 1,3-...	11.48	35.8	ug/l	887220	4	12.08	1239430	50.0
Cyclohexane, butyl-	12.00	34.4	ug/l	852887	4	12.08	1239430	50.0
Benzene, 1-methyl...	12.32	78.0	ug/l	1932430	4	12.08	1239430	50.0
o-Cymene	12.61	60.6	ug/l	1502300	4	12.08	1239430	50.0
Benzene, 2-ethyl...	12.66	41.0	ug/l	1016610	4	12.08	1239430	50.0
4-Methylphenyl ac...	12.77	39.2	ug/l	971051	4	12.08	1239430	50.0