

Data Path : Z:\VOASRV\HPCHEM1\MSVOA W\DATA\VW091118\
 Data File : VW005297.D
 Acq On : 12 Sep 2018 02:32
 Operator : SY/AP
 Sample : J4821-04
 Misc : 10.42G/5ML/MSVOA W/SOIL
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 ENY-SB-102-61.0-61.5

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_W\METHOD\82W091018S.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	2.221	45	52	53	rBV4	3956	7345	4.65%	0.604%
2	4.928	493	496	501	rVB4	1442	2450	1.55%	0.201%
3	7.641	934	941	946	rBV	3756	8228	5.21%	0.677%
4	7.891	973	982	987	rBV	23945	57216	36.23%	4.705%
5	7.958	987	993	1002	rVB	59145	130438	82.60%	10.727%
6	8.311	1044	1051	1062	rVB2	26530	58297	36.91%	4.794%
7	8.848	1126	1139	1151	rBV	74742	155364	98.38%	12.777%
8	10.329	1375	1382	1389	rBV	70547	126124	79.86%	10.372%
9	10.384	1389	1391	1400	rVB7	1202	2187	1.38%	0.180%
10	10.482	1400	1407	1418	rBV	62401	127962	81.03%	10.523%
11	10.713	1439	1445	1451	rBV	27631	68783	43.55%	5.656%
12	10.866	1467	1470	1476	rBV7	1522	2348	1.49%	0.193%
13	11.085	1500	1506	1507	rBV6	2319	3335	2.11%	0.274%
14	11.146	1509	1516	1517	rBV7	2995	4632	2.93%	0.381%
15	11.640	1588	1597	1607	rBV	85369	157924	100.00%	12.987%
16	11.725	1609	1611	1614	rVB3	1987	2031	1.29%	0.167%
17	11.841	1627	1630	1632	rBV4	1529	1696	1.07%	0.139%
18	12.628	1752	1759	1765	rBV2	58260	103268	65.39%	8.492%
19	12.762	1777	1781	1786	rBV6	4273	8533	5.40%	0.702%
20	12.865	1796	1798	1803	rVB5	2090	2051	1.30%	0.169%
21	12.945	1809	1811	1813	rBV3	3009	2519	1.60%	0.207%
22	13.255	1860	1862	1868	rVB5	5590	7134	4.52%	0.587%
23	13.505	1899	1903	1907	rBV7	3730	7553	4.78%	0.621%
24	13.566	1907	1913	1920	rVB	54921	93226	59.03%	7.667%
25	13.829	1954	1956	1959	rVB4	2998	3219	2.04%	0.265%
26	13.883	1961	1965	1970	rBV8	4614	7007	4.44%	0.576%
27	14.085	1994	1998	2010	rVB2	10302	27980	17.72%	2.301%
28	14.213	2013	2019	2024	rBV10	3285	6975	4.42%	0.574%
29	14.280	2028	2030	2032	rBV3	1512	1714	1.09%	0.141%
30	14.396	2045	2049	2054	rBV8	4413	8604	5.45%	0.708%
31	14.505	2065	2067	2069	rBV3	1808	1588	1.01%	0.131%
32	14.700	2096	2099	2100	rBV3	1819	2234	1.41%	0.184%
33	14.804	2110	2116	2125	rVB10	5546	12395	7.85%	1.019%
34	15.420	2215	2217	2220	rVB4	1790	1889	1.20%	0.155%

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35	15.511	2229	2232	2233	rBV3	1508	1753	1.11%	0.144%
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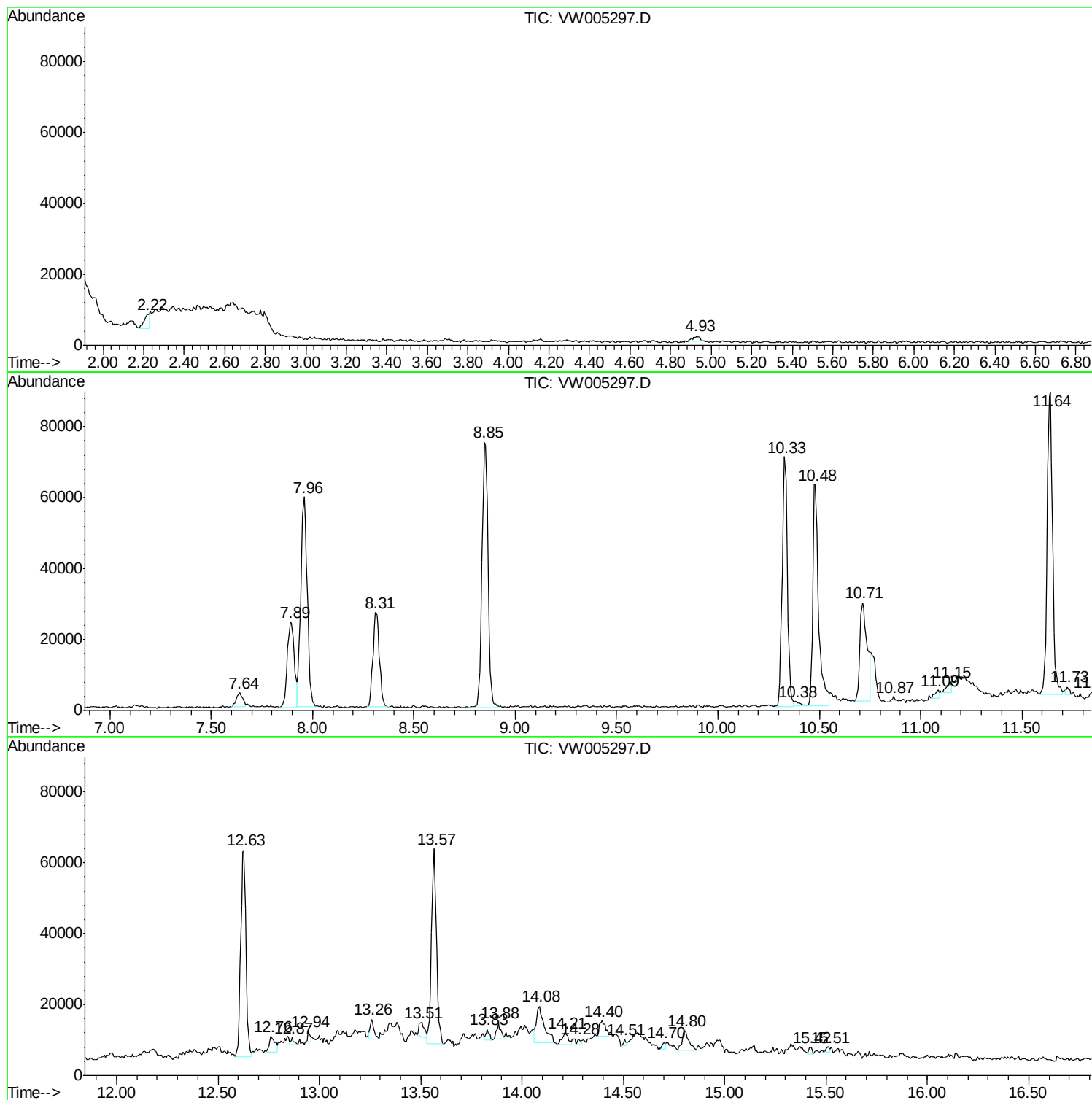
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W091018S.M
Quant Title : SW846 8260

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



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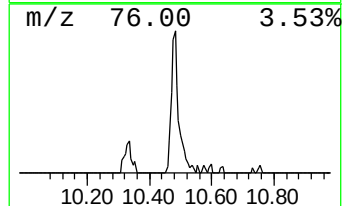
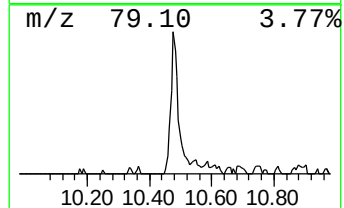
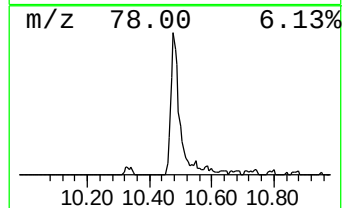
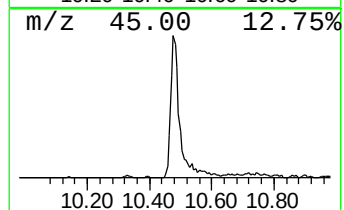
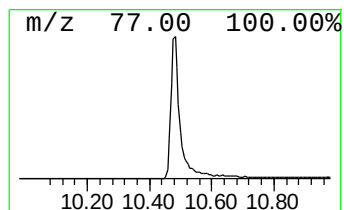
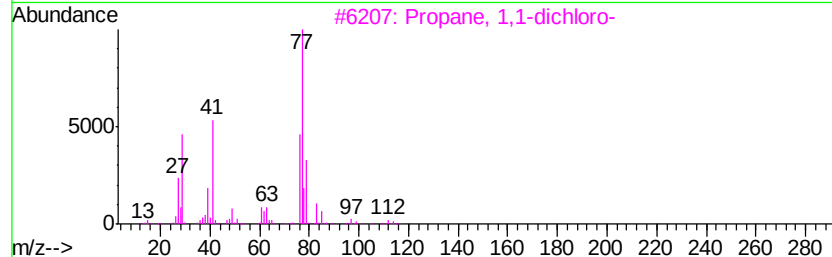
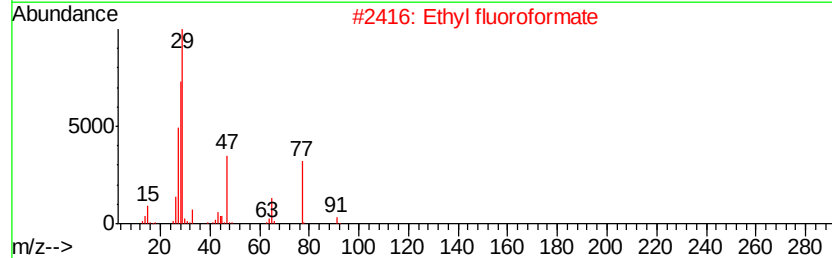
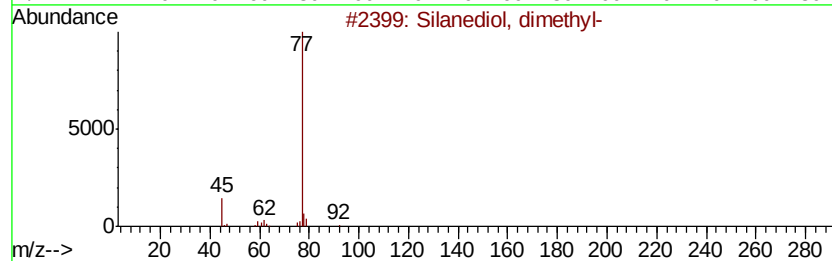
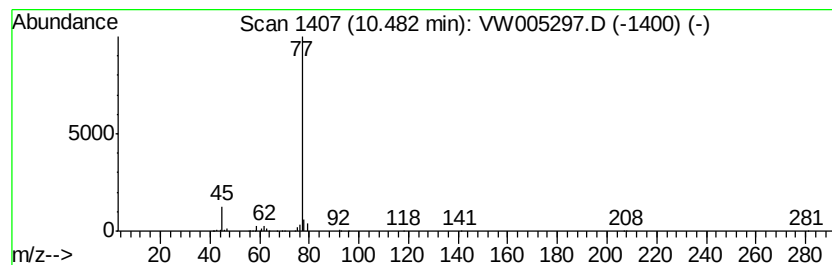
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_W\METHOD\82W091018S.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown10.48 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.48	40.51 ug/l	127962	Chlorobenzene-d5	11.64

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Silanediol, dimethyl-	92	C2H8O2Si	001066-42-8	49
2		Ethyl fluoroformate	92	C3H5FO2	000461-64-3	2
3		Propane, 1,1-dichloro-	112	C3H6Cl2	000078-99-9	1
4		Dimethoxamine	77	C2H7NO2	007487-32-3	1
5		1,5-Hexadiene, 3,3,4,4-tetrafluoro-	154	C6H6F4	001763-21-9	1



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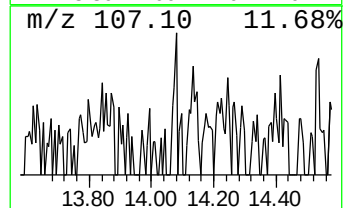
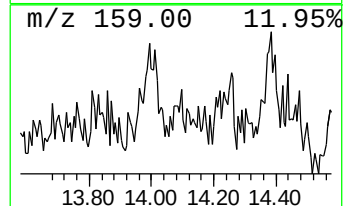
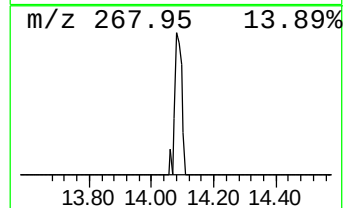
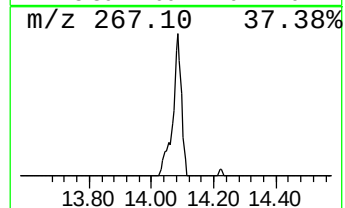
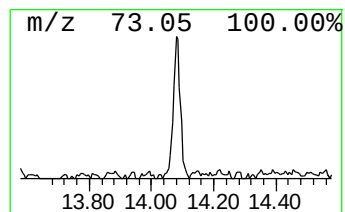
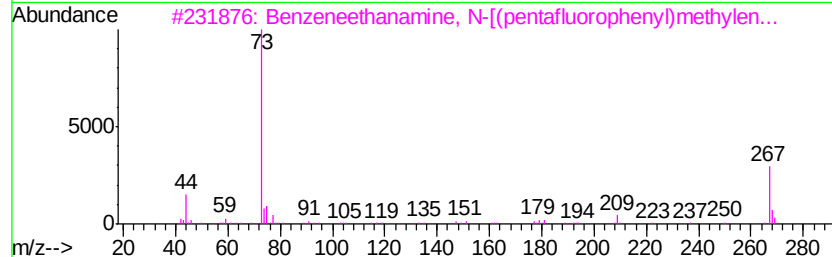
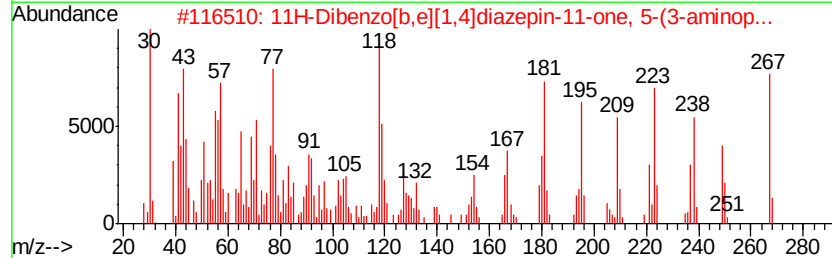
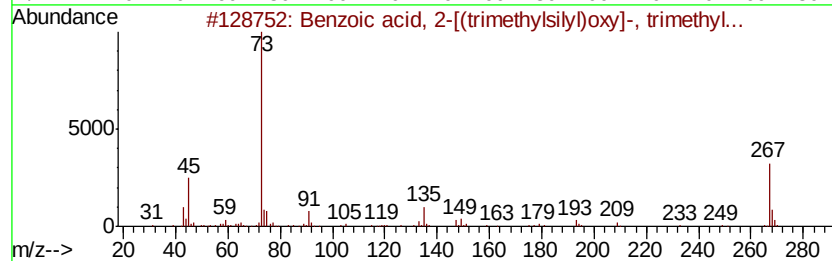
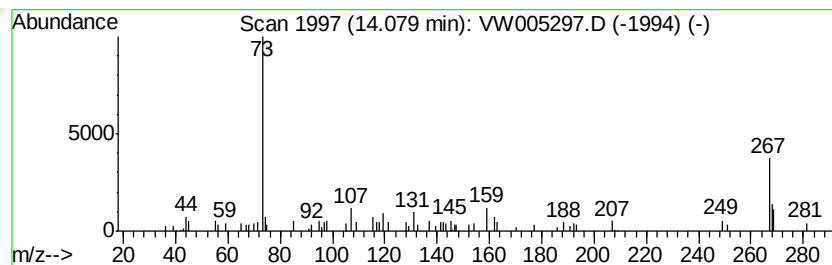
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Quant Title : SW846 8260

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

Peak Number 3 unknown14.08 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.08	15.01 ug/l	27980	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzoic acid, 2-[(trimethylsilyl)oxy]-, trimethyl...	282	C13H22O3Si2	003789-85-3	38
2		11H-Dibenzo[b,e][1,4]diazepin-11-one, 5-(3-aminopropyl)-	267	C16H17N3O	013450-73-2	27
3		Benzeneethanamine, N-[(pentafluorophenyl)methyl]-	475	C21H26F5N02Si2	055429-85-1	9
4		4-Methylcatechol, bis(trimethylsilyl) ether	268	C13H24O2Si2	1000332-79-9	7
5		Benzeneacetic acid, trimethylsilyl ester	208	C11H16O2Si	002078-18-4	7



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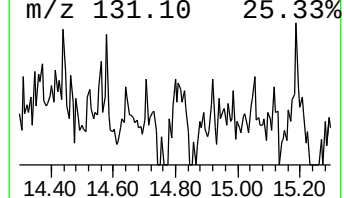
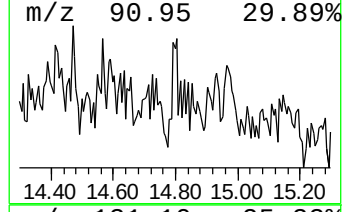
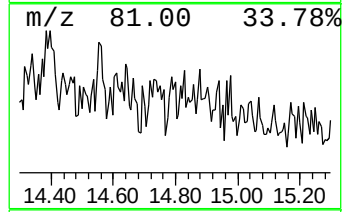
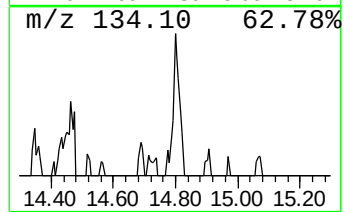
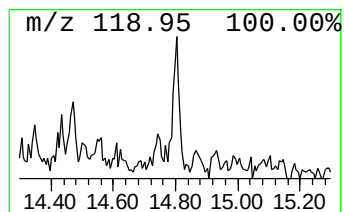
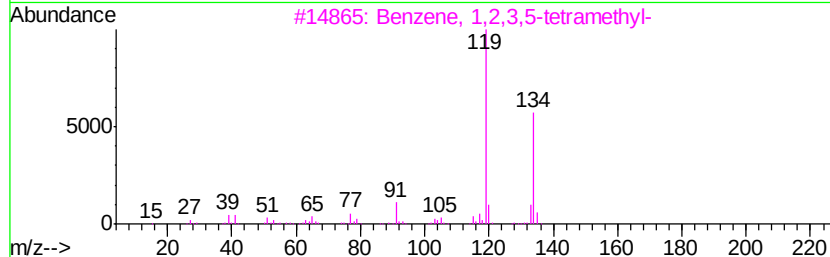
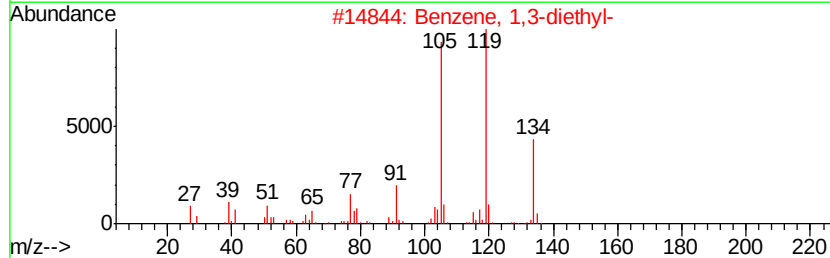
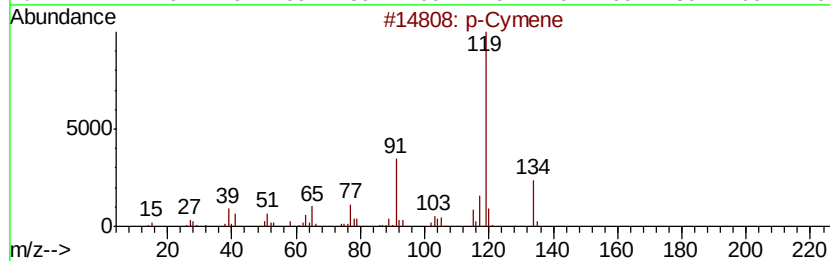
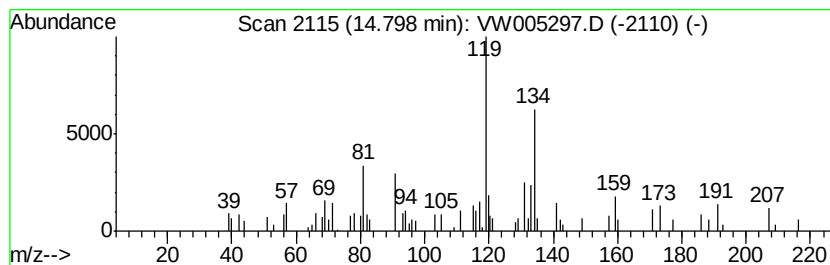
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Peak Number 4 unknown14.80 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.80	6.65 ug/l	12395	1,4-Dichlorobenzene-d4	13.57

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	p-Cymene	134	C10H14	000099-87-6	49
2		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	46
3		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	46
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	46
5		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	46



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TIC Top Hit name	RT	EstConc	Units	Response	---Internal Standard---			
					#	RT	Resp	Conc
unknown10.48	10.48	40.5	ug/l	127962	3	11.64	157924	50.0
unknown14.08	14.08	15.0	ug/l	27980	4	13.57	93226	50.0
unknown14.80	14.80	6.7	ug/l	12395	4	13.57	93226	50.0