

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD090920\
 Data File : VD066581.D
 Acq On : 09 Sep 2020 14:19
 Operator : VA/SY
 Sample : L3877-11
 Misc : 5.84G/5.00ml/MSVOA D/SOIL
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 TP-3

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_D\METHOD\82D090320S.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	7.473	935	944	956	rBV2	136532	367334	15.30%	2.620%
2	7.914	1008	1019	1024	rBV	348514	839644	34.98%	5.989%
3	7.979	1024	1030	1044	rVB	577067	1309572	54.55%	9.341%
4	8.326	1079	1089	1100	rBV	297819	687801	28.65%	4.906%
5	8.861	1169	1180	1193	rBV	868886	1754706	73.10%	12.516%
6	9.343	1256	1262	1270	rVB6	13163	29058	1.21%	0.207%
7	10.337	1422	1431	1448	rBV	1316442	2400467	100.00%	17.121%
8	11.243	1580	1585	1590	rVB5	28295	46561	1.94%	0.332%
9	11.349	1598	1603	1608	rBV7	14631	25493	1.06%	0.182%
10	11.526	1628	1633	1637	rBV3	16126	26585	1.11%	0.190%
11	11.643	1644	1653	1665	rBV	1210309	2078179	86.57%	14.823%
12	11.743	1665	1670	1675	rBV3	37438	67722	2.82%	0.483%
13	11.855	1679	1689	1694	rBV8	18843	58214	2.43%	0.415%
14	12.179	1732	1744	1751	rBV9	24492	87751	3.66%	0.626%
15	12.631	1814	1821	1832	rBV	886946	1515956	63.15%	10.813%
16	12.802	1845	1850	1857	rVB8	20359	50091	2.09%	0.357%
17	12.967	1873	1878	1883	rBV9	21293	34230	1.43%	0.244%
18	13.126	1898	1905	1913	rVB3	31014	63300	2.64%	0.451%
19	13.273	1925	1930	1935	rBV3	17680	31853	1.33%	0.227%
20	13.414	1946	1954	1956	rBV9	14379	34798	1.45%	0.248%
21	13.520	1967	1972	1974	rBV4	15206	24240	1.01%	0.173%
22	13.578	1974	1982	1992	rVB	1109665	1850430	77.09%	13.198%
23	13.778	2011	2016	2021	rBV2	75938	129900	5.41%	0.927%
24	13.896	2032	2036	2041	rVB6	16843	29217	1.22%	0.208%
25	13.973	2041	2049	2053	rBV10	11979	30409	1.27%	0.217%
26	14.120	2071	2074	2078	rVB5	17882	25298	1.05%	0.180%
27	14.231	2087	2093	2097	rBV8	18922	37756	1.57%	0.269%
28	14.443	2125	2129	2132	rBV5	20377	30843	1.28%	0.220%
29	14.484	2133	2136	2140	rVB2	27747	38285	1.59%	0.273%
30	14.720	2170	2176	2179	rBV7	15375	28270	1.18%	0.202%
31	14.820	2189	2193	2194	rBV	28097	31571	1.32%	0.225%
32	14.896	2200	2206	2214	rVB9	28434	59149	2.46%	0.422%
33	14.990	2217	2222	2225	rBV7	17325	31733	1.32%	0.226%
34	15.096	2236	2240	2246	rBV8	18203	32670	1.36%	0.233%

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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

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35	16.220	2424	2431	2434	rBV8	10698	26681	1.11%	0.190%
36	16.708	2507	2514	2525	rVB5	41944	104439	4.35%	0.745%

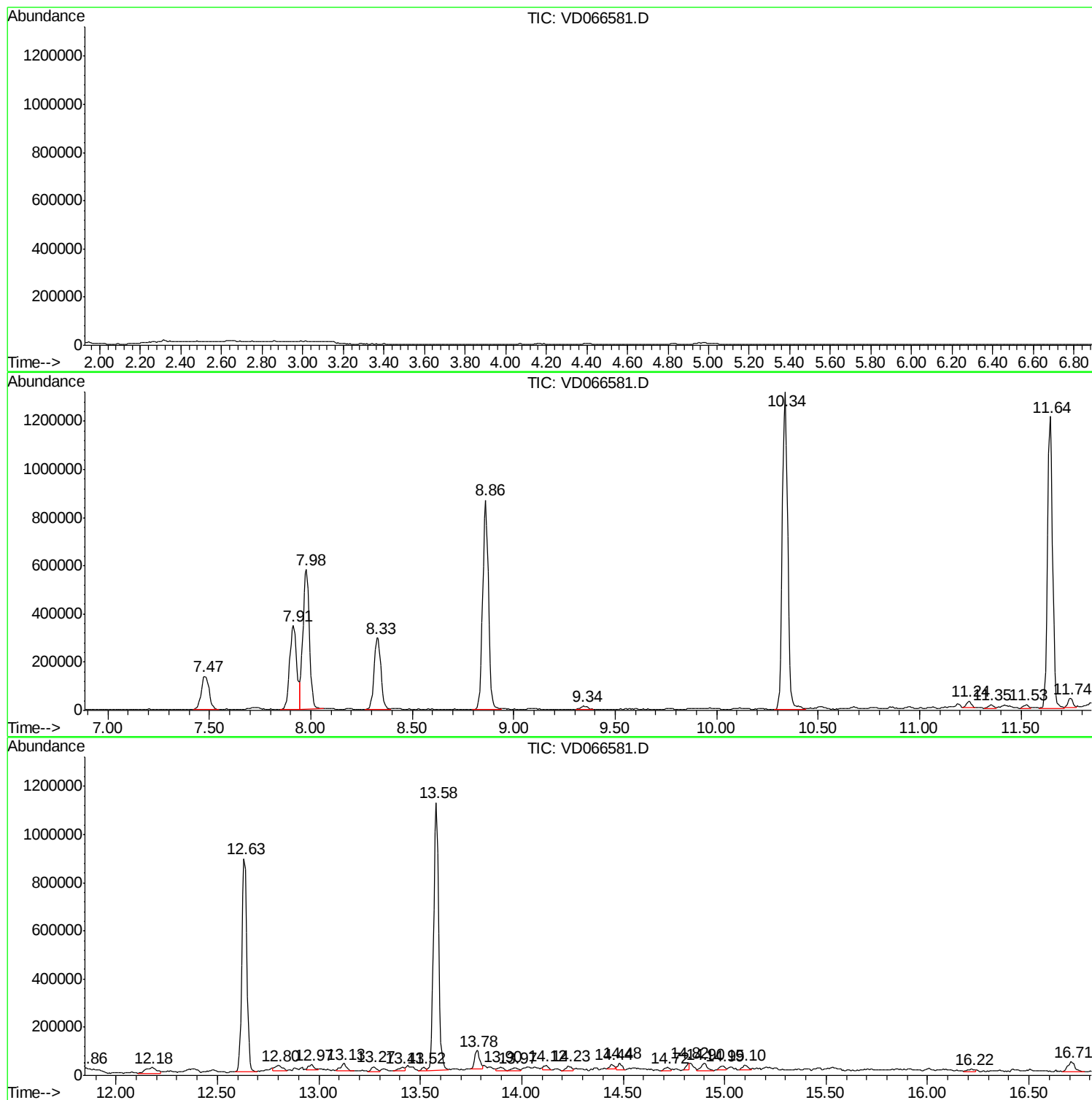
Sum of corrected areas: 14020206

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D090320S.M
Quant Title : SW846 8260

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



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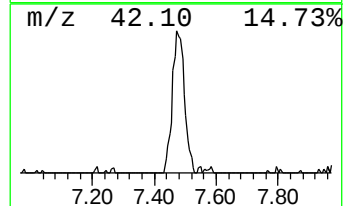
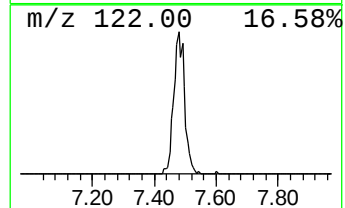
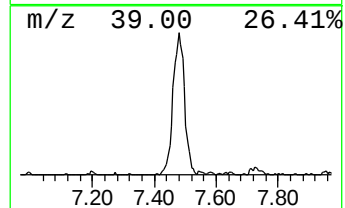
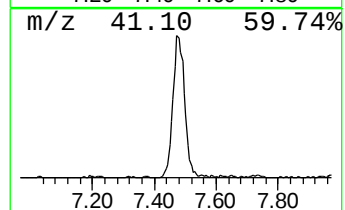
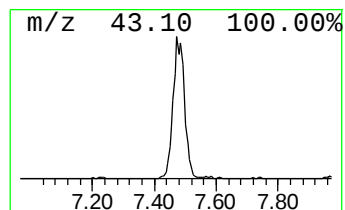
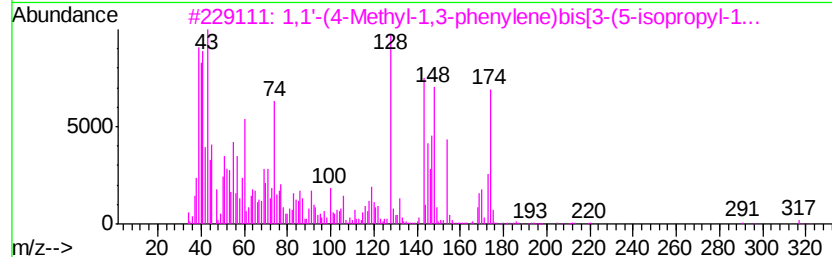
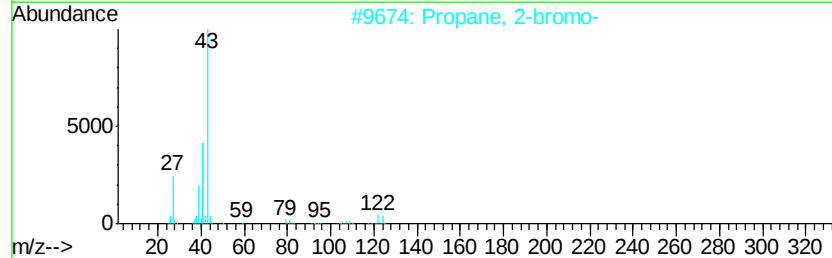
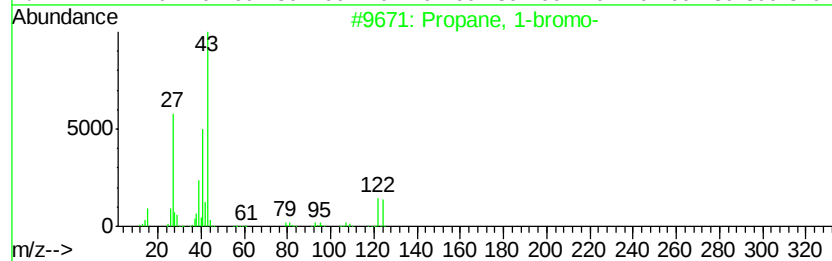
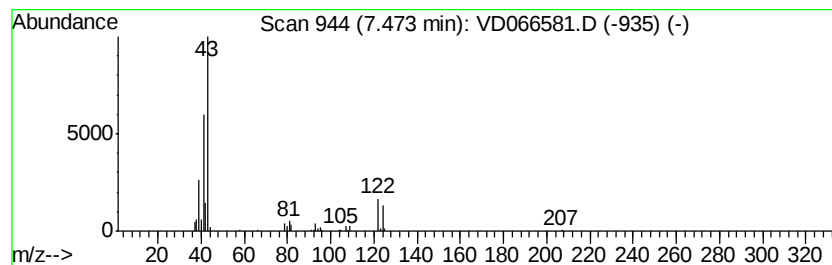
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Peak Number 1 Propane, 1-bromo- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.47	14.03 ug/l	367334	Pentafluorobenzene	7.98

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane, 1-bromo-	122	C3H7Br	000106-94-5	90
2		Propane, 2-bromo-	122	C3H7Br	000075-26-3	59
3		1,1'-(4-Methyl-1,3-phenylene)bis...	460	C19H24N8O2S2	1000224-58-4	10
4		Propene	42	C3H6	000115-07-1	9
5		Propane, 2,2-dibromo-	200	C3H6Br2	000594-16-1	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propane, 1-bromo-	7.47	14.0	ug/l	367334	1	7.98	1309570	50.0