

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY042121\
 Data File : VY004550.D
 Acq On : 21 Apr 2021 13:03
 Operator : SY/MD
 Sample : M2103-01
 Misc : 6.72G/5ML/MSVOA_Y/SOIL
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 RO-124030

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_Y\methods\82Y041221S.M

Title : SW846 8260

Signal : TIC: VY004550.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.186	47	60	61	rBV4	7051	23028	1.30%	0.371%
2	7.728	960	969	975	rBV2	100903	240849	13.58%	3.880%
3	7.801	975	981	991	rVB	173835	388253	21.89%	6.254%
4	8.149	1029	1038	1049	rVB	111263	245192	13.82%	3.950%
5	8.697	1120	1128	1141	rVB	282733	548793	30.94%	8.840%
6	10.185	1364	1372	1379	rBV	468403	783272	44.16%	12.617%
7	10.807	1468	1474	1483	rVB	16670	28126	1.59%	0.453%
8	10.904	1483	1490	1498	rBV	15604	33135	1.87%	0.534%
9	11.093	1515	1521	1527	rBV2	25710	46761	2.64%	0.753%
10	11.203	1534	1539	1544	rBV2	19884	35493	2.00%	0.572%
11	11.301	1551	1555	1562	rVB2	14118	24932	1.41%	0.402%
12	11.496	1579	1587	1597	rBV	421709	686853	38.72%	11.064%
13	11.679	1612	1617	1622	rVV7	8713	20862	1.18%	0.336%
14	11.740	1622	1627	1631	rVV	19653	37332	2.10%	0.601%
15	11.782	1631	1634	1639	rVB2	13057	20118	1.13%	0.324%
16	12.008	1663	1671	1677	rBV3	26284	61372	3.46%	0.989%
17	12.124	1687	1690	1694	rVB	26085	34402	1.94%	0.554%
18	12.209	1694	1704	1705	rBV6	19715	40535	2.29%	0.653%
19	12.239	1705	1709	1714	rVB2	30629	57425	3.24%	0.925%
20	12.374	1725	1731	1737	rVB5	16408	32050	1.81%	0.516%
21	12.489	1743	1750	1758	rBV3	1013791	1773701	100.00%	28.571%
22	12.648	1772	1776	1783	rVB3	25087	44087	2.49%	0.710%
23	12.727	1783	1789	1795	rBV7	13028	34089	1.92%	0.549%
24	12.813	1795	1803	1808	rVV4	30512	80749	4.55%	1.301%
25	12.861	1808	1811	1816	rVB2	15080	25350	1.43%	0.408%
26	13.044	1837	1841	1849	rVB3	18186	33582	1.89%	0.541%
27	13.325	1879	1887	1890	rBV4	19854	38245	2.16%	0.616%
28	13.428	1897	1904	1912	rVB	401357	604742	34.09%	9.741%
29	13.751	1949	1957	1961	rBV4	24182	45401	2.56%	0.731%
30	13.965	1986	1992	1998	rVB2	60481	104648	5.90%	1.686%
31	14.263	2035	2041	2051	rVB7	15060	34593	1.95%	0.557%

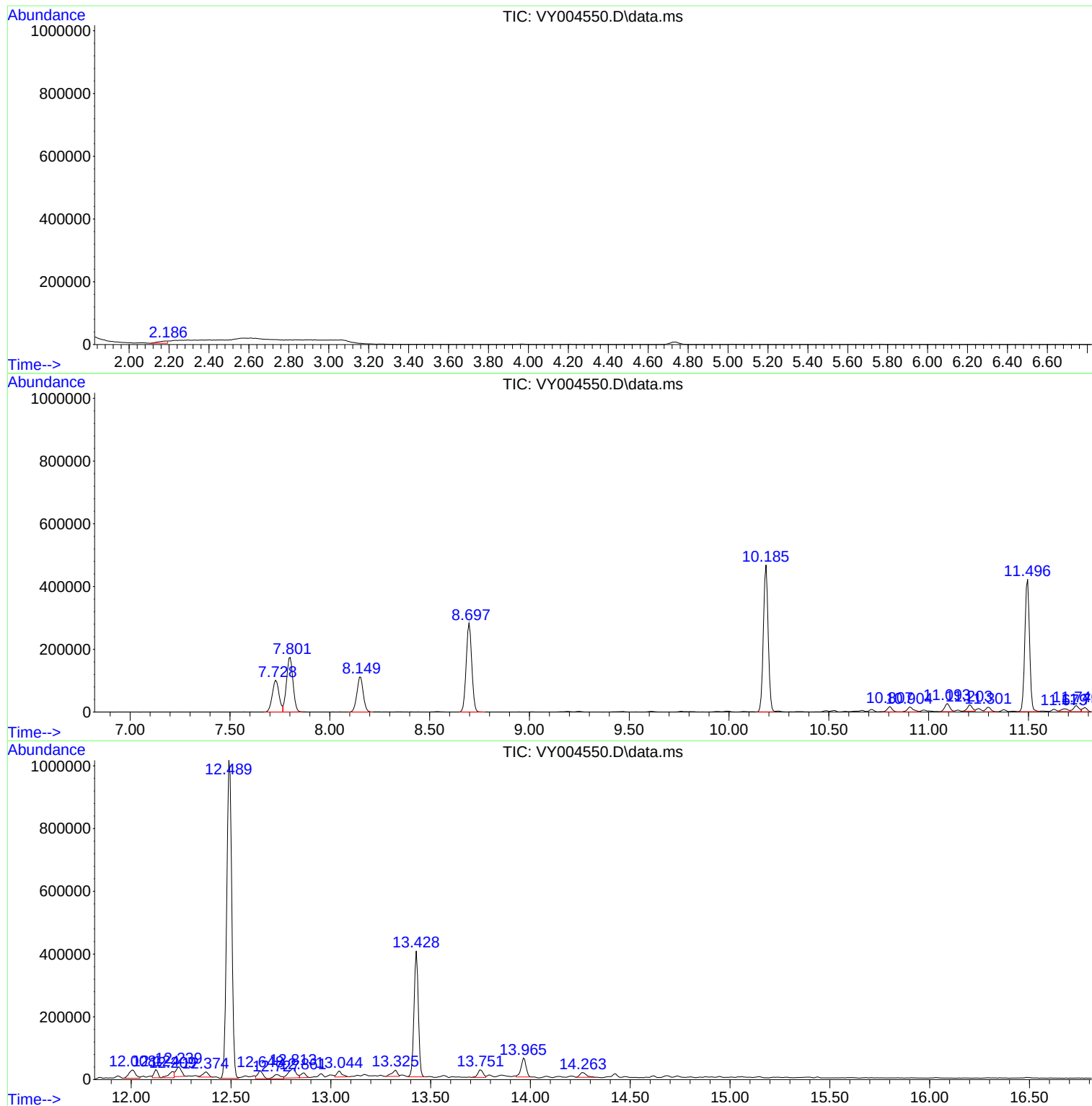
Sum of corrected areas: 6207970

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Instrument :
MSVOA_Y
ClientSampleId :
RO-124030

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041221S.M
Quant Title : SW846 8260

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



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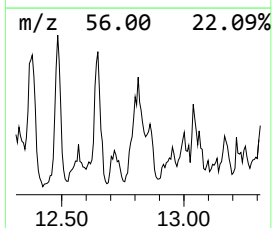
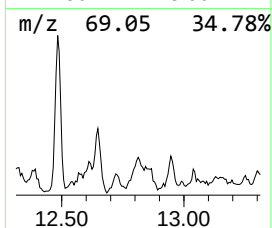
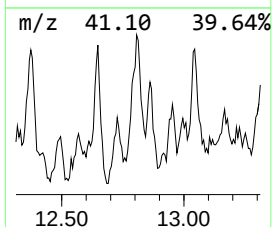
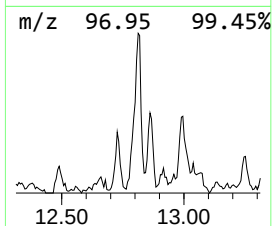
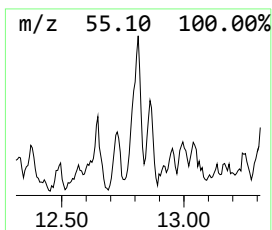
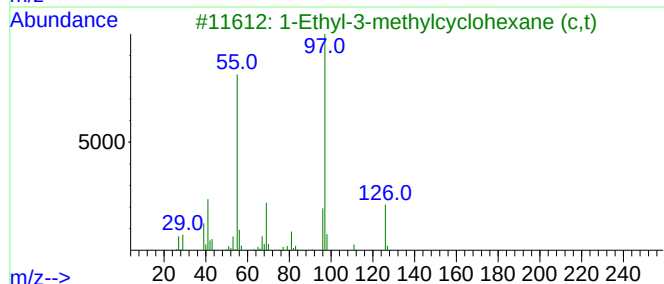
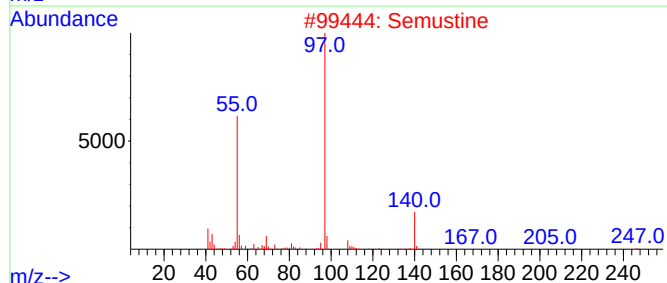
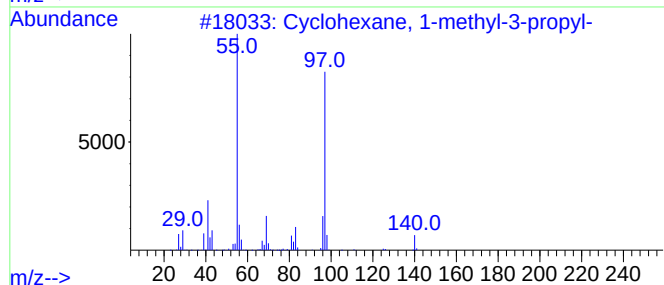
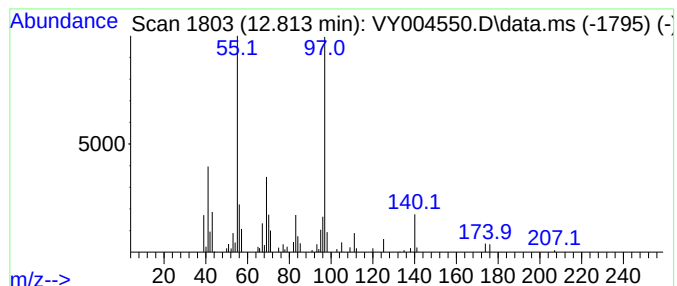
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y041221S.M
Quant Title : SW846 8260

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TIC Integration Parameters: LSCINT.P

Peak Number 1 Cyclohexane, 1-methyl-3-pro... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.813	6.68 ug/l	80749	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-methyl-3-propyl-	140	C10H20	004291-80-9	64
2			Semustine	247	C10H18ClN3O2	013909-09-6	59
3			1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	59
4			Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	59
5			Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	006069-98-3	58



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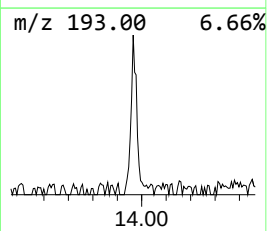
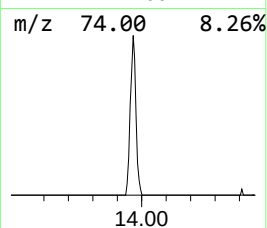
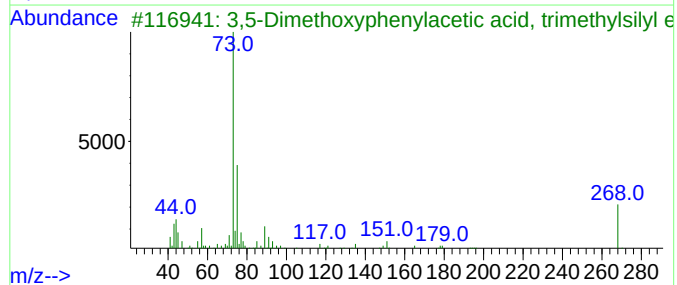
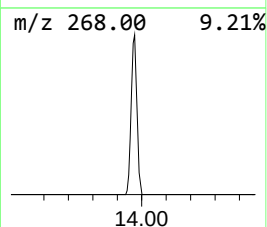
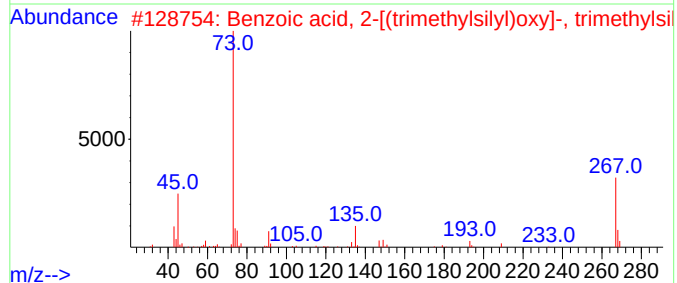
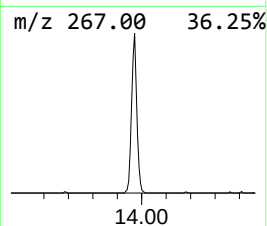
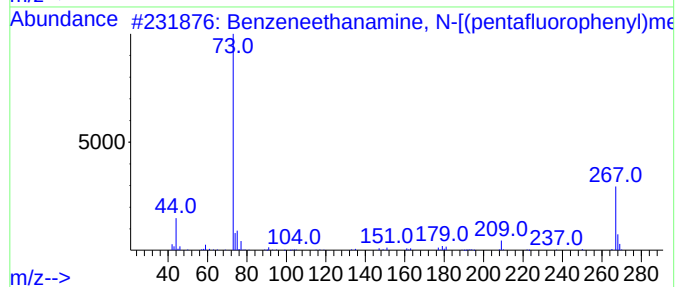
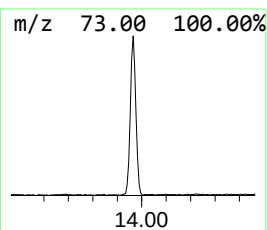
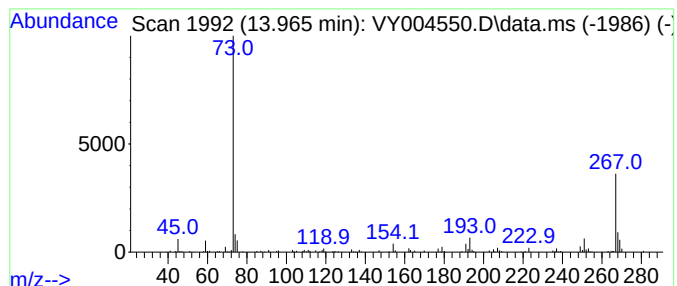
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Quant Title : SW846 8260

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Peak Number 2 Benzeneethanamine, N-[(pent... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.965	8.65 ug/l	104648	1,4-Dichlorobenzene-d4	13.428

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzeneethanamine, N-[(pentafluoro...	475	C21H26F5NO2Si2	055429-85-1	59
2			Benzoic acid, 2-[(trimethylsilyl)...	282	C13H22O3Si2	003789-85-3	50
3			3,5-Dimethoxyphenylacetic acid, ...	268	C13H20O4Si	1000071-82-4	35
4			11H-Dibenzo[b,e][1,4]diazepin-11...	267	C16H17N3O	013450-73-2	30
5			1-Boraindane, 3-methyl-1-[1-(tri...	266	C17H23BSi	190316-36-0	27



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Cyclohexane, 1-...	12.813	6.7	ug/l	80749	4	13.428	604742	50.0
Benzeneethanami...	13.965	8.7	ug/l	104648	4	13.428	604742	50.0