

Data Path : Z:\voasrv\HPCHEM1\MSVOA_Y\Data\VY061622\
 Data File : VY009203.D
 Acq On : 16 Jun 2022 15:28
 Operator : KP/MD
 Sample : N3367-01
 Misc : 5.00g/5.0mL/MSVOA_Y/SOIL
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 MJS-LB-02-13.0-14.0

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\82Y061522S.M
 Title : SW846 8260

Signal : TIC: VY009203.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.198	380	390	396	rBV5	1445	4214	1.66%	0.237%
2	4.698	464	472	473	rBV2	2828	4192	1.65%	0.236%
3	5.740	633	643	652	rVB5	3031	9309	3.66%	0.524%
4	6.972	834	845	856	rBV	13193	41394	16.26%	2.331%
5	7.716	957	967	972	rBV	36450	83648	32.86%	4.710%
6	7.783	972	978	988	rVB	64889	148322	58.26%	8.352%
7	8.136	1024	1036	1045	rBV2	41078	117711	46.24%	6.628%
8	8.685	1113	1126	1134	rBV	99575	193664	76.07%	10.905%
9	10.179	1363	1371	1379	rBV	144501	254584	100.00%	14.335%
10	10.581	1431	1437	1442	rBV2	8071	14213	5.58%	0.800%
11	10.941	1491	1496	1507	rBV3	6166	11920	4.68%	0.671%
12	11.313	1552	1557	1565	rVB2	4999	8646	3.40%	0.487%
13	11.489	1579	1586	1596	rVB	126708	210305	82.61%	11.842%
14	12.477	1741	1748	1758	rVB2	91650	152526	59.91%	8.588%
15	12.642	1769	1775	1779	rBV5	1641	2663	1.05%	0.150%
16	12.806	1798	1802	1806	rVV5	3545	6529	2.56%	0.368%
17	13.117	1849	1853	1858	rBV6	1890	3488	1.37%	0.196%
18	13.306	1876	1884	1885	rBV4	2112	3618	1.42%	0.204%
19	13.422	1896	1903	1912	rVB	111558	175638	68.99%	9.890%
20	13.629	1932	1937	1941	rBV5	2019	3365	1.32%	0.189%
21	13.751	1950	1957	1960	rBV6	4669	8758	3.44%	0.493%
22	13.965	1988	1992	1997	rVB6	2620	3565	1.40%	0.201%
23	14.074	2007	2010	2017	rVB6	1648	2990	1.17%	0.168%
24	14.141	2017	2021	2025	rBV6	1789	2701	1.06%	0.152%
25	14.251	2035	2039	2047	rVB9	3214	7551	2.97%	0.425%
26	14.422	2062	2067	2070	rVB6	2067	3133	1.23%	0.176%
27	14.477	2070	2076	2080	rVB6	1553	2633	1.03%	0.148%
28	14.611	2094	2098	2102	rVB6	3577	6142	2.41%	0.346%
29	14.678	2102	2109	2113	rBV7	5011	10948	4.30%	0.616%
30	14.733	2113	2118	2125	rVB5	4228	7912	3.11%	0.446%
31	14.946	2146	2153	2158	rVB8	3146	8142	3.20%	0.458%
32	15.233	2190	2200	2208	rBV	115273	200149	78.62%	11.270%
33	15.318	2208	2214	2226	rVV3	9481	22810	8.96%	1.284%
34	15.434	2229	2233	2239	rVB9	2181	3715	1.46%	0.209%
35	16.293	2367	2374	2380	rBV4	10517	22057	8.66%	1.242%
36	16.489	2399	2406	2413	rVB5	6137	12808	5.03%	0.721%

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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\82Y061522S.M
Title : SW846 8260

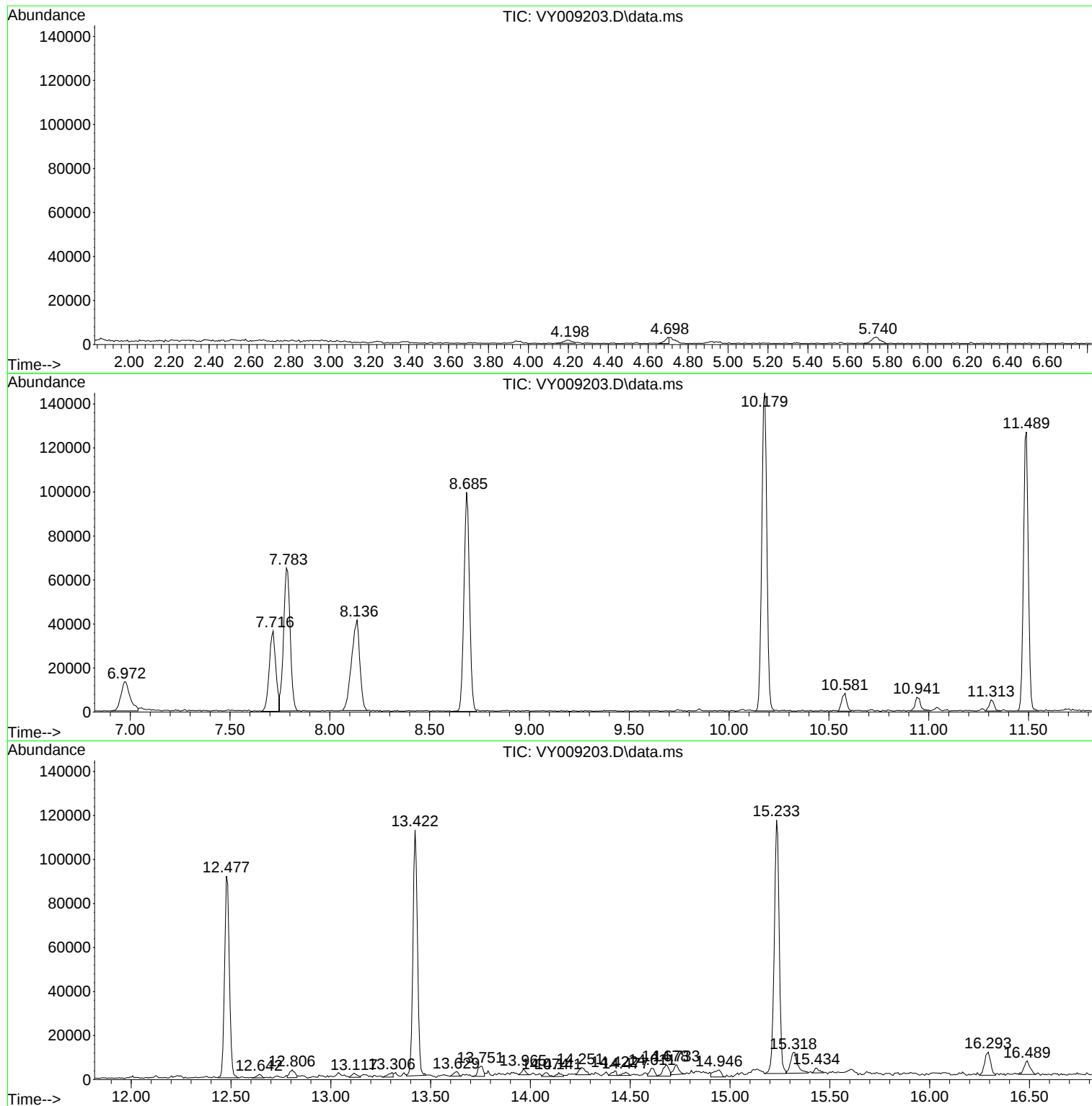
Sum of corrected areas: 1775963

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_Y\methods\82Y061522S.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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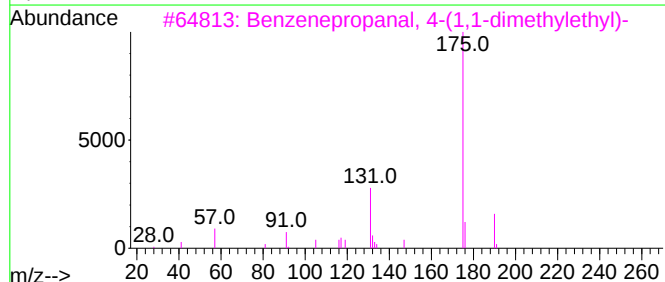
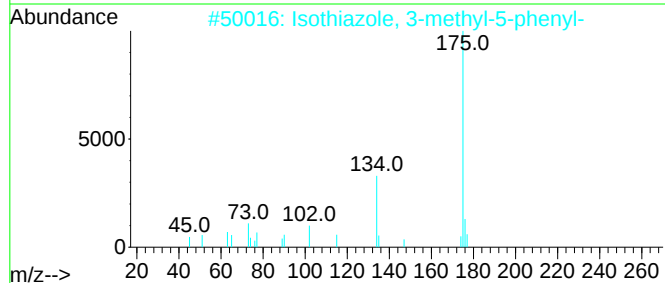
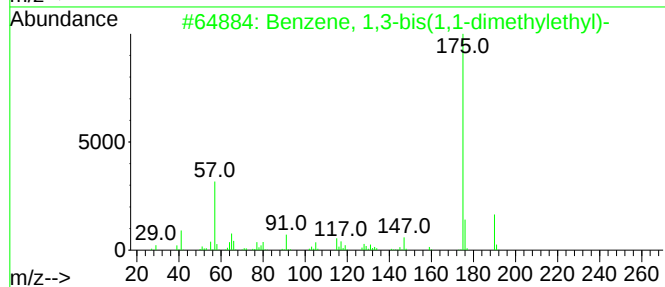
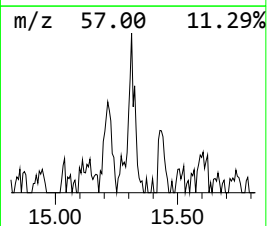
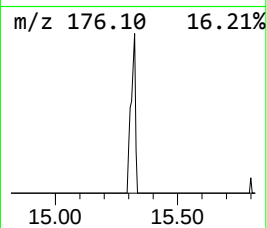
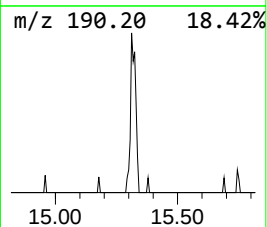
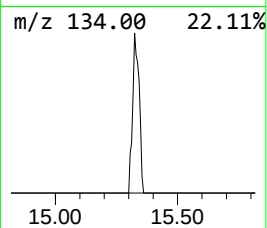
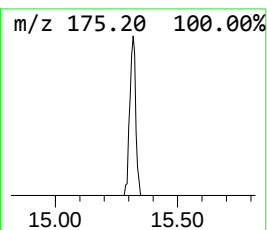
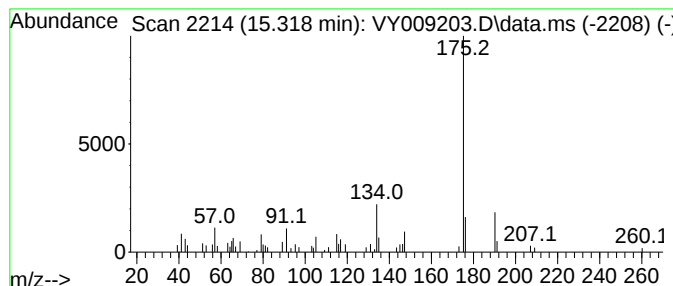
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Peak Number 3 Benzene, 1,3-bis(1,1-dimeth... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.318	6.49 ug/l	22810	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-bis(1,1-dimethyleth...	190	C14H22	001014-60-4	76
2			Isothiazole, 3-methyl-5-phenyl-	175	C10H9NS	001732-45-2	64
3			Benzenepropanal, 4-(1,1-dimethyl...	190	C13H18O	018127-01-0	62
4			2,2'-Ethylidenebis(5-methylfuran)	190	C12H14O2	003209-79-8	59
5			Benzene, 1,5-dimethyl-2,4-bis(1-...	190	C14H22	005186-68-5	59



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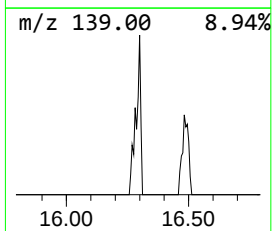
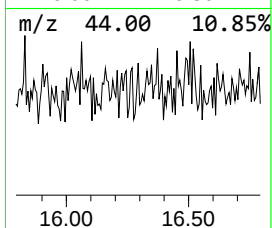
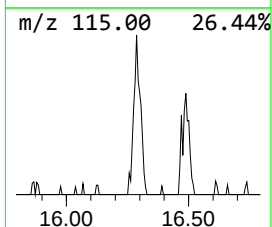
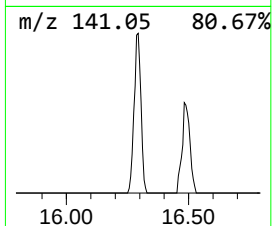
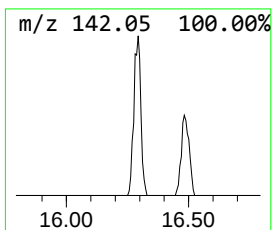
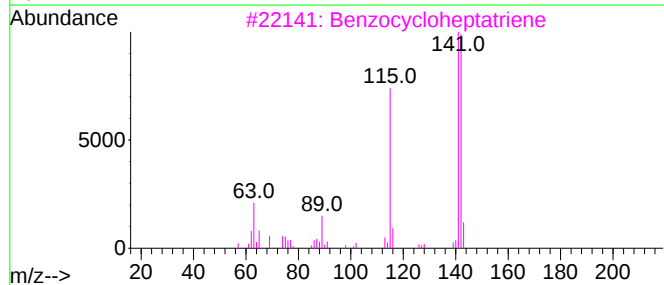
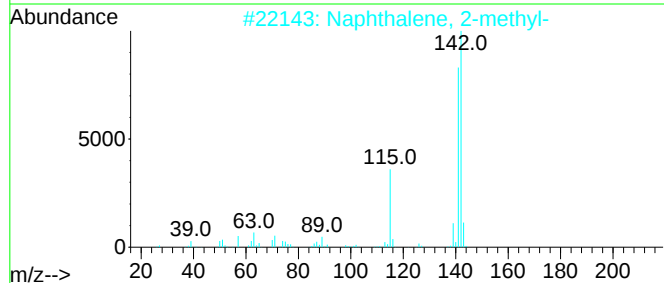
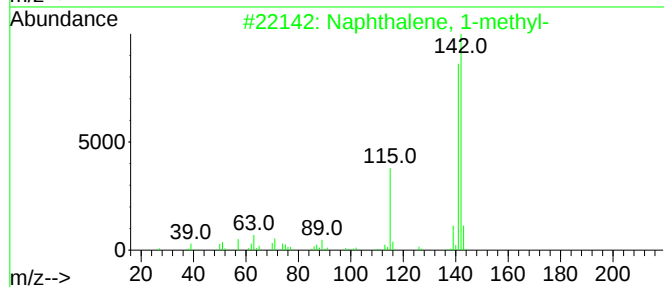
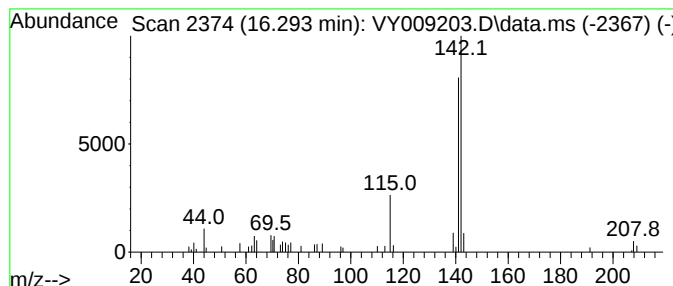
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Peak Number 4 Benzocycloheptatriene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.294	6.28 ug/l	22057	1,4-Dichlorobenzene-d4	13.422

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	91
3			Benzocycloheptatriene	142	C11H10	000264-09-5	87
4			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	78
5			3-Indolecarbonitrile	142	C9H6N2	005457-28-3	46



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1,3-bi...	15.318	6.5	ug/l	22810	4	13.422	175638	50.0
Benzocyclohepta...	16.294	6.3	ug/l	22057	4	13.422	175638	50.0