LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041222\
Data File : VU047953.D
Acq On : 12 Apr 2022 14:17
Operator : SY/MD
Sample : N2293-08MEDL 5X
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ETNN8MEDL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0 % 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_U\Method\SFAMULM032422WMA.M
Title : VOC Analysis

Signal : TIC: VU047953.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.600	74	81	97	rBV	106128	135652	1.63%	0.602%
2	1.899	167	174	192	rVB	95513	138907	1.66%	0.616%
3	2.565	369	381	396	rBV	209850	348542	4.18%	1.546%
4	2.957	497	503	514	rVB2	1792	2787	0.03%	0.012%
5	3.044	523	530	542	rBV6	2336	5023	0.06%	0.022%
6	3.443	650	654	656	rBV2	1022	916	0.01%	0.004%
7	3.514	674	676	685	rVV3	547	476	0.01%	0.002%
8	3.768	749	755	759	rBV4	368	408	0.00%	0.002%
9	3.825	767	773	774	rBV3	903	763	0.01%	0.003%
10	4.070	846	849	854	rVV3	586	622	0.01%	0.003%
11	4.169	874	880	885	rBV5	633	564	0.01%	0.003%
12	4.391	941	949	954	rVB3	533	887	0.01%	0.004%
13	4.626	1008	1022	1060	rBV	132637	330466	3.96%	1.466%
14	5.063	1141	1158	1181	rBV	209962	457281	5.48%	2.028%
15	5.491	1287	1291	1296	rVB5	607	525	0.01%	0.002%
16	5.726	1339	1364	1376	rBV2	412239	1140547	13.67%	5.059%
17	5.896	1405	1417	1430	rBV5	6277	13643	0.16%	0.061%
18	6.066	1468	1470	1475	rVB3	728	428	0.01%	0.002%
19	6.112	1479	1484	1487	rBV2	490	494	0.01%	0.002%
20	6.163	1496	1500	1505	rBV4	406	428	0.01%	0.002%
21	6.250	1511	1527	1553	rBV	468108	875707	10.50%	3.884%
22	6.475	1588	1597	1607	rBV5	3397	5769	0.07%	0.026%
23	6.690	1649	1664	1684	rVV	265205	519569	6.23%	2.304%
24	6.809	1696	1701	1707	rVB5	1149	1579	0.02%	0.007%
25	6.851	1707	1714	1717	rBV3	479	438	0.01%	0.002%
26	6.935	1735	1740	1746	rBV3	368	494	0.01%	0.002%
27	7.185	1807	1818	1828	rBV5	4239	8416	0.10%	0.037%
28	7.491	1904	1913	1924	rBV4	7305	12515	0.15%	0.056%
29	7.568	1924	1937	1960	rBV	142163	254831	3.05%	1.130%
30	7.687	1968	1974	1982	rBV7	1836	2781	0.03%	0.012%
31	7.800	2006	2009	2013	rBV3	433	402	0.00%	0.002%
32	7.899	2024	2040	2054	rBV	500281	856931	10.27%	3.801%
33	7.976	2056	2064	2081	rVB3	32564	68611	0.82%	0.304%
34	8.179	2114	2127	2142	rBV	102443	172167	2.06%	0.764%
35	8.359	2173	2183	2195	rVB5	8444	15519	0.19%	0.069%
36	8.545	2238	2241	2244	rVV2	746	491	0.01%	0.002%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041222\
Data File : VU047953.D
Acq On : 12 Apr 2022 14:17
Operator : SY/MD
Sample : N2293-08MEDL 5X
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : ETNN8MEDL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0 % 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_U\Method\SFAMULM032422WMA.M
Title : VOC Analysis

37	8.636	2255	2269	2302	rBV	457058	808132	9.69%	3.584%
38	8.918	2354	2357	2359	rBV4	579	398	0.00%	0.002%
39	9.018	2379	2388	2400	rVB6	7086	12095	0.14%	0.054%
40	9.362	2487	2495	2500	rBV3	5873	8347	0.10%	0.037%
41	9.420	2500	2513	2515	rBV	732652	995417	11.93%	4.415%
42	9.578	2555	2562	2570	rBV5	9703	13640	0.16%	0.060%
43	9.693	2593	2598	2610	rVB2	5494	8639	0.10%	0.038%
44	9.861	2647	2650	2656	rBV6	919	887	0.01%	0.004%
45	9.889	2656	2659	2668	rVB6	1402	1764	0.02%	0.008%
46	9.989	2685	2690	2692	rBV3	2555	2556	0.03%	0.011%
47	10.095	2709	2723	2730	rBV3	5447	12710	0.15%	0.056%
48	10.221	2752	2762	2768	rBV7	5453	10240	0.12%	0.045%
49	10.330	2793	2796	2800	rBV4	1094	1123	0.01%	0.005%
50	10.404	2809	2819	2830	rBV8	3136	6419	0.08%	0.028%
51	10.478	2840	2842	2848	rVB5	814	663	0.01%	0.003%
52	10.558	2857	2867	2869	rBV5	1120	1443	0.02%	0.006%
53	10.658	2887	2898	2902	rBV5	4714	7845	0.09%	0.035%
54	10.700	2903	2911	2918	rVV2	23030	40049	0.48%	0.178%
55	10.758	2918	2929	2947	rVB	474026	774837	9.29%	3.437%
56	10.893	2963	2971	2973	rBV7	1688	1788	0.02%	0.008%
57	10.979	2990	2998	3001	rBV6	2447	3806	0.05%	0.017%
58	11.098	3028	3035	3038	rBV4	1711	2110	0.03%	0.009%
59	11.378	3115	3122	3130	rBB8	1383	1948	0.02%	0.009%
60	11.417	3130	3134	3139	rVB4	782	596	0.01%	0.003%
61	11.471	3141	3151	3157	rBV5	4832	8262	0.10%	0.037%
62	11.578	3177	3184	3194	rBV9	1803	2916	0.03%	0.013%
63	11.626	3196	3199	3204	rVB3	682	493	0.01%	0.002%
64	11.680	3212	3216	3217	rBV2	675	461	0.01%	0.002%
65	11.693	3219	3220	3225	rVB3	699	582	0.01%	0.003%
66	11.719	3225	3228	3230	rVB2	823	560	0.01%	0.002%
67	11.745	3230	3236	3244	rBV4	2639	3222	0.04%	0.014%
68	11.812	3244	3257	3274	rBV	856192	1369833	16.42%	6.076%
69	11.951	3292	3300	3313	rVB5	20749	35230	0.42%	0.156%
70	12.066	3333	3336	3340	rBV5	780	568	0.01%	0.003%
71	12.102	3340	3347	3355	rVB9	1126	1647	0.02%	0.007%
72	12.195	3361	3376	3407	rBV	626989	1135404	13.61%	5.036%
73	12.323	3410	3416	3425	rVB5	4151	6051	0.07%	0.027%
74	12.381	3430	3434	3435	rBV3	815	550	0.01%	0.002%
75	12.468	3454	3461	3467	rBV8	1949	3307	0.04%	0.015%
76	12.577	3487	3495	3498	rBV5	1073	1585	0.02%	0.007%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041222\
Data File : VU047953.D
Acq On : 12 Apr 2022 14:17
Operator : SY/MD
Sample : N2293-08MEDL 5X
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ETNN8MEDL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0 % 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_U\Method\SFAMULM032422WMA.M
Title : VOC Analysis

77	12.629	3501	3511	3521	rVV9	5218	9111	0.11%	0.040%
78	12.703	3531	3534	3536	rBV3	1405	885	0.01%	0.004%
79	12.770	3548	3555	3557	rBV5	1557	1762	0.02%	0.008%
80	12.835	3568	3575	3585	rVB6	6860	10779	0.13%	0.048%
81	12.967	3603	3616	3631	rVB	210489	357196	4.28%	1.584%
82	13.076	3647	3650	3653	rBV4	821	414	0.00%	0.002%
83	13.269	3704	3710	3717	rBV9	2935	4372	0.05%	0.019%
84	13.327	3717	3728	3736	rBV	167025	270873	3.25%	1.201%
85	13.375	3736	3743	3755	rVB2	134838	203101	2.43%	0.901%
86	13.950	3915	3922	3923	rBV4	2110	2036	0.02%	0.009%
87	14.044	3944	3951	3959	rBV8	8716	13047	0.16%	0.058%
88	14.214	4002	4004	4006	rBV3	802	443	0.01%	0.002%
89	14.359	4036	4049	4064	rBV	134066	210575	2.52%	0.934%
90	14.494	4085	4091	4099	rBV9	3785	4877	0.06%	0.022%
91	14.590	4113	4121	4133	rVB4	7653	12581	0.15%	0.056%
92	14.883	4207	4212	4221	rVB9	2937	4156	0.05%	0.018%
93	15.111	4276	4283	4292	rBV8	9068	14494	0.17%	0.064%
94	15.310	4329	4345	4355	rBV2	133222	231122	2.77%	1.025%
95	15.545	4405	4418	4437	rBV	832801	1287851	15.44%	5.712%
96	15.796	4487	4496	4507	rBV6	34972	57931	0.69%	0.257%
97	15.892	4510	4526	4541	rBV3	193335	532553	6.38%	2.362%
98	16.388	4669	4680	4692	rBV	5499784	8343279	100.00%	37.006%
99	17.208	4928	4935	4942	rBV3	34410	52273	0.63%	0.232%
100	17.259	4942	4951	4966	rVB2	156256	256502	3.07%	1.138%

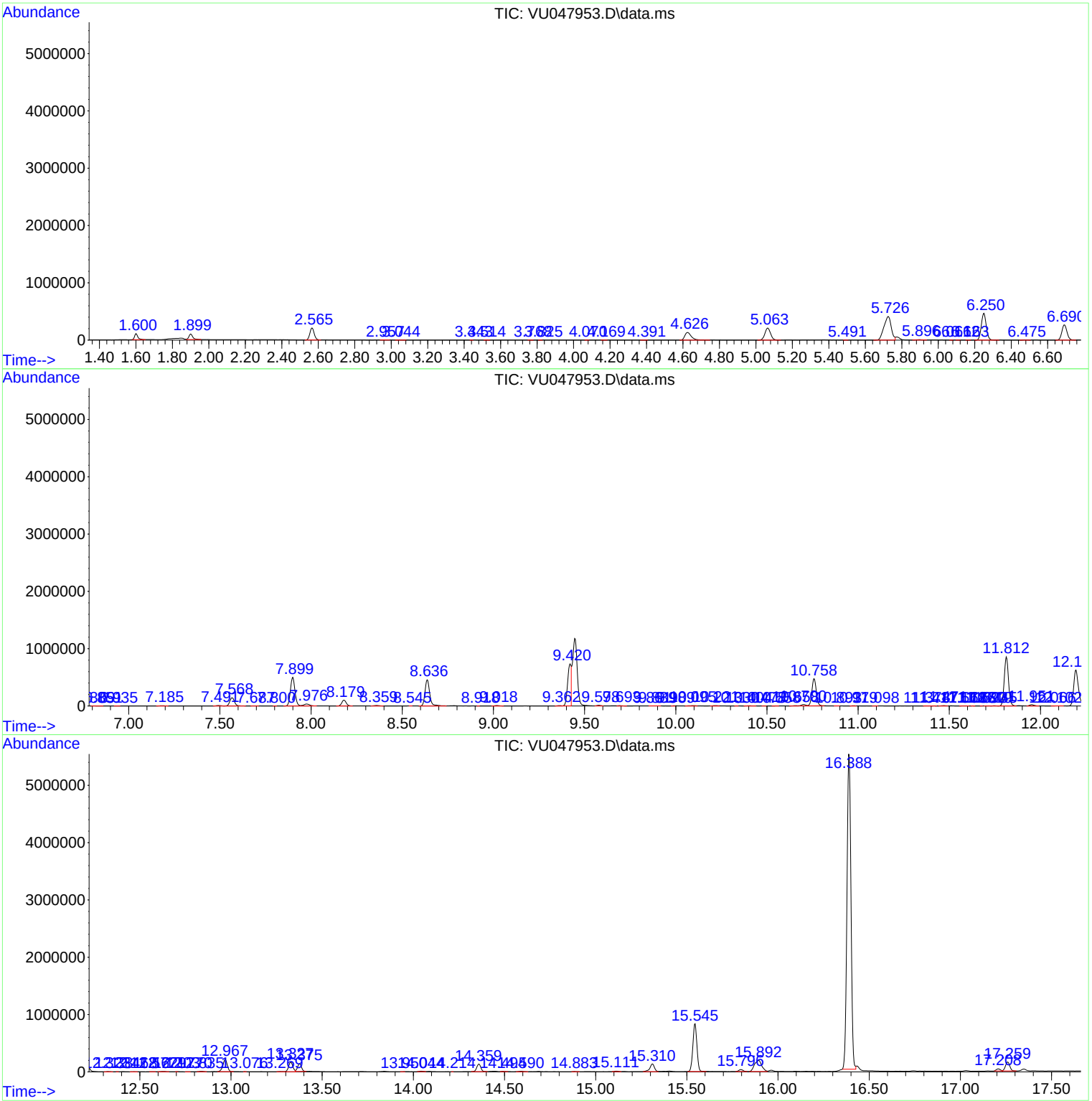
Sum of corrected areas: 22545915

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041222\
Data File : VU047953.D
Acq On : 12 Apr 2022 14:17
Operator : SY/MD
Sample : N2293-08MEDL 5X
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ETNN8MEDL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM032422WMA.M
Quant Title : VOC Analysis

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041222\
Data File : VU047953.D
Acq On : 12 Apr 2022 14:17
Operator : SY/MD
Sample : N2293-08MEDL 5X
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ETNN8MEDL

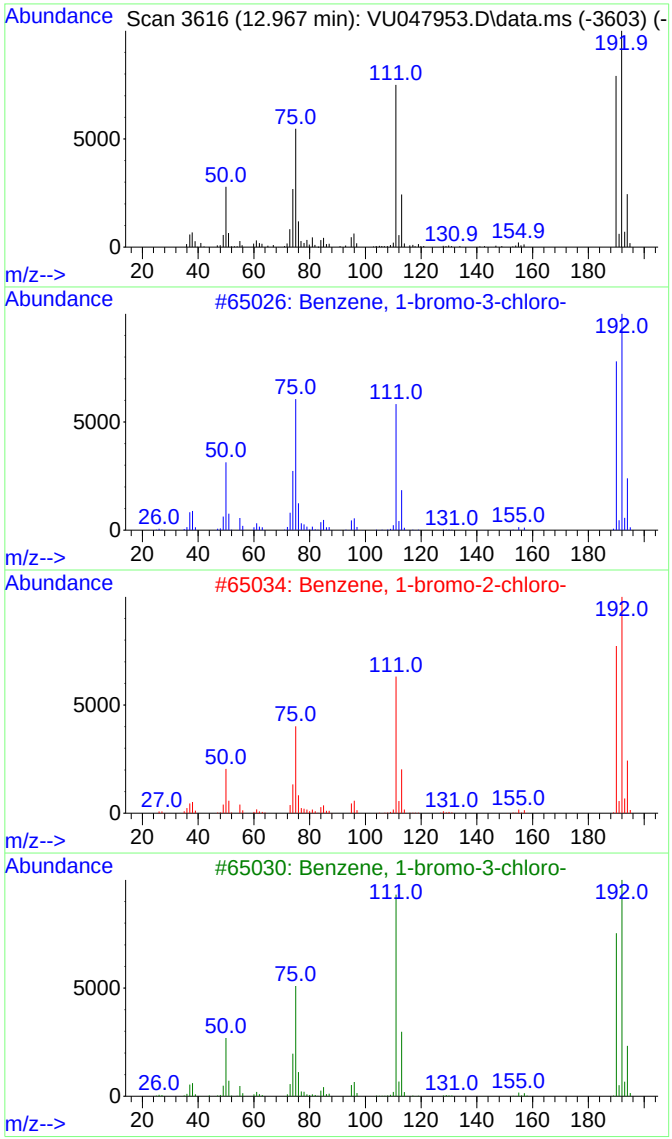
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM032422WMA.M
Quant Title : VOC Analysis

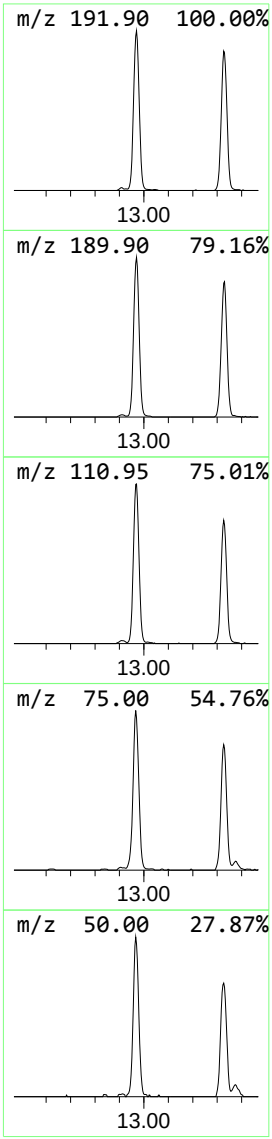
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Benzene, 1-bromo-3-chloro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.967	13.04 ug/L	357196	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	99
2	Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
3	Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	98
4	Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
5	Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041222\
Data File : VU047953.D
Acq On : 12 Apr 2022 14:17
Operator : SY/MD
Sample : N2293-08MEDL 5X
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ETNN8MEDL

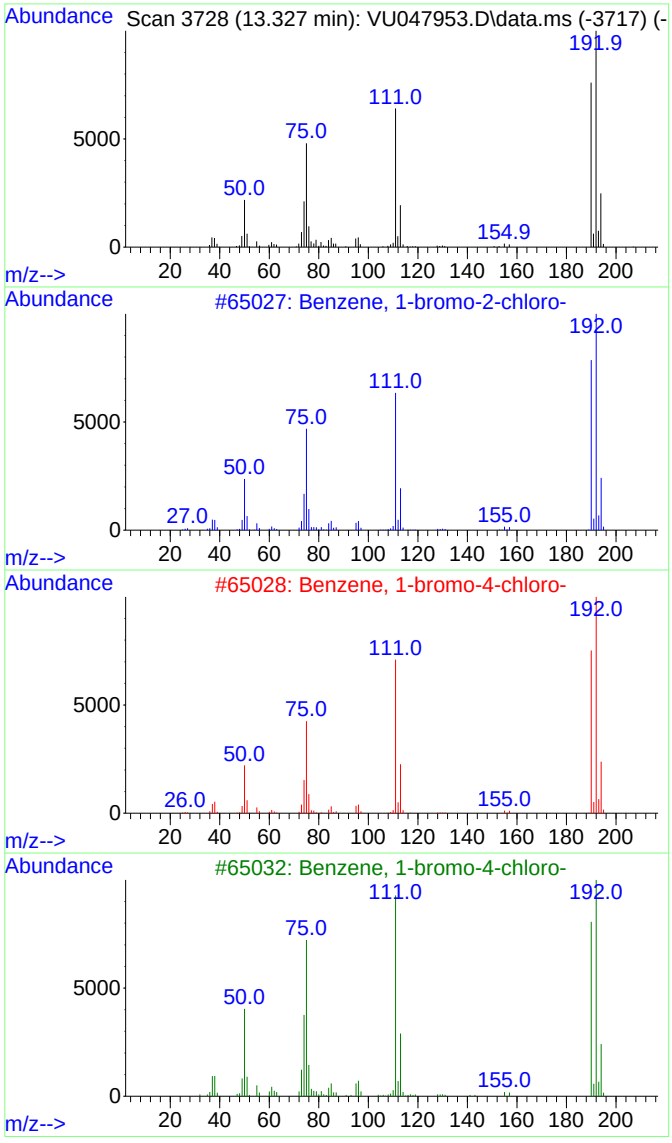
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM032422WMA.M
Quant Title : VOC Analysis

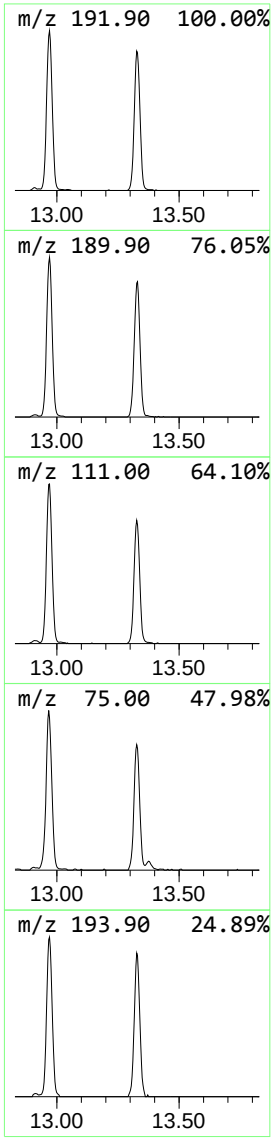
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1-bromo-2-chloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.327	9.89 ug/L	270873	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	99
2	Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
3	Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
4	Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
5	4-Chlorophenylselenol	192	C6H5ClSe	016645-10-6	97





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041222\
Data File : VU047953.D
Acq On : 12 Apr 2022 14:17
Operator : SY/MD
Sample : N2293-08MEDL 5X
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ETNN8MEDL

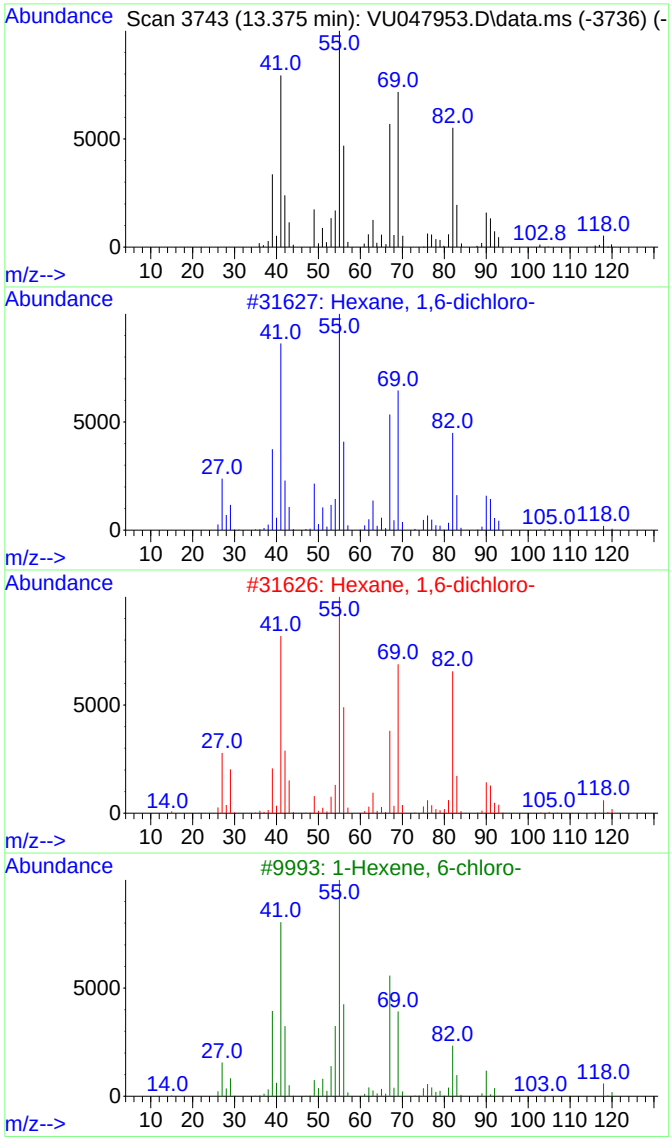
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM032422WMA.M
Quant Title : VOC Analysis

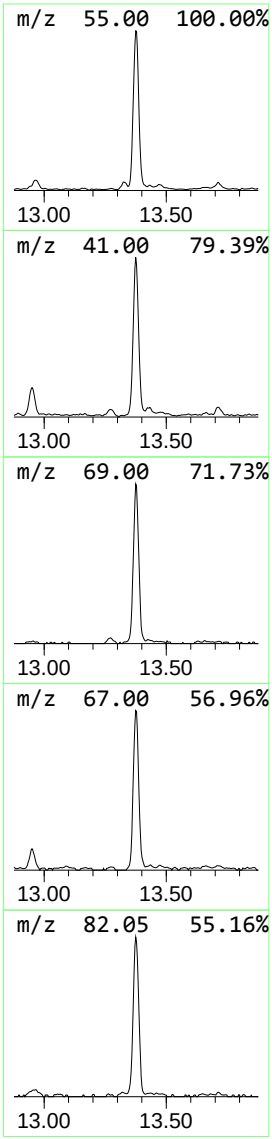
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Hexane, 1,6-dichloro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.375	7.41 ug/L	203101	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 1,6-dichloro-	154	C6H12Cl2	002163-00-0	86
2	Hexane, 1,6-dichloro-	154	C6H12Cl2	002163-00-0	80
3	1-Hexene, 6-chloro-	118	C6H11Cl	000928-89-2	58
4	Cyclohexane, chloro-	118	C6H11Cl	000542-18-7	50
5	Hexenyl angelate, 4z-	182	C11H18O2	1000383-63-7	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041222\
Data File : VU047953.D
Acq On : 12 Apr 2022 14:17
Operator : SY/MD
Sample : N2293-08MEDL 5X
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ETNN8MEDL

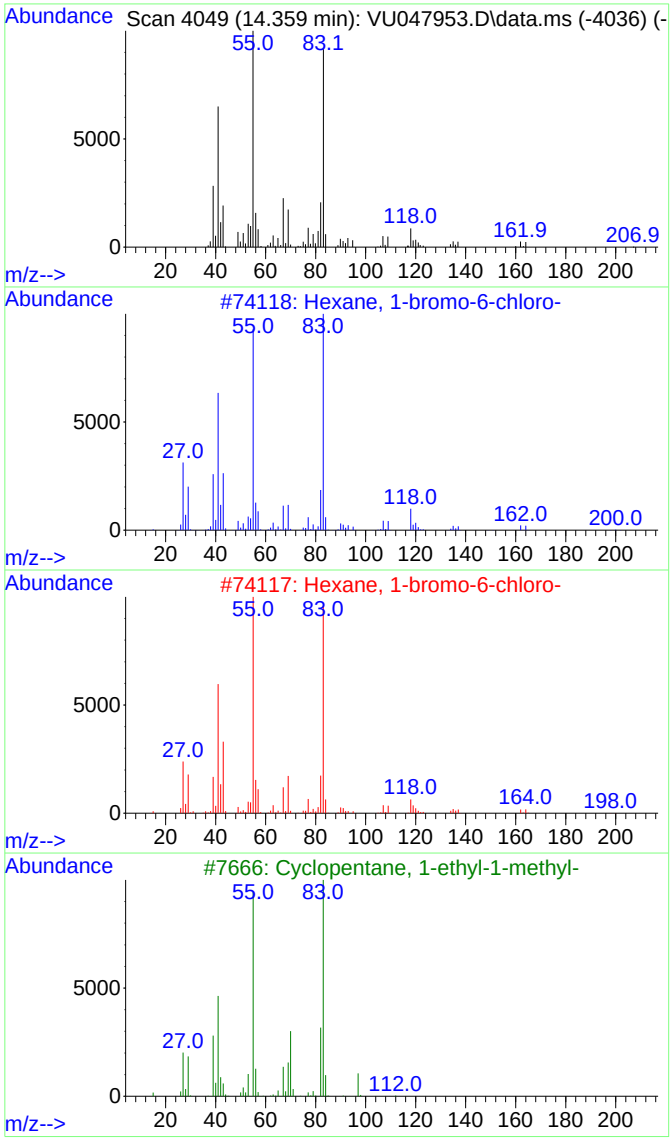
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM032422WMA.M
Quant Title : VOC Analysis

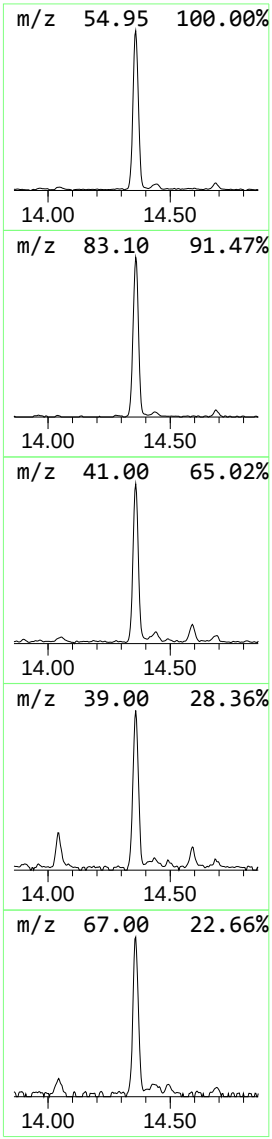
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Hexane, 1-bromo-6-chloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.359	7.69 ug/L	210575	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 1-bromo-6-chloro-	198	C6H12BrCl	006294-17-3	86
2	Hexane, 1-bromo-6-chloro-	198	C6H12BrCl	006294-17-3	72
3	Cyclopentane, 1-ethyl-1-methyl-	112	C8H16	016747-50-5	64
4	Cyclopentane, 1-ethyl-1-methyl-	112	C8H16	016747-50-5	59
5	Oxalic acid, cyclohexyl isohexyl...	256	C14H24O4	1010309-30-7	53





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041222\
Data File : VU047953.D
Acq On : 12 Apr 2022 14:17
Operator : SY/MD
Sample : N2293-08MEDL 5X
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ETNN8MEDL

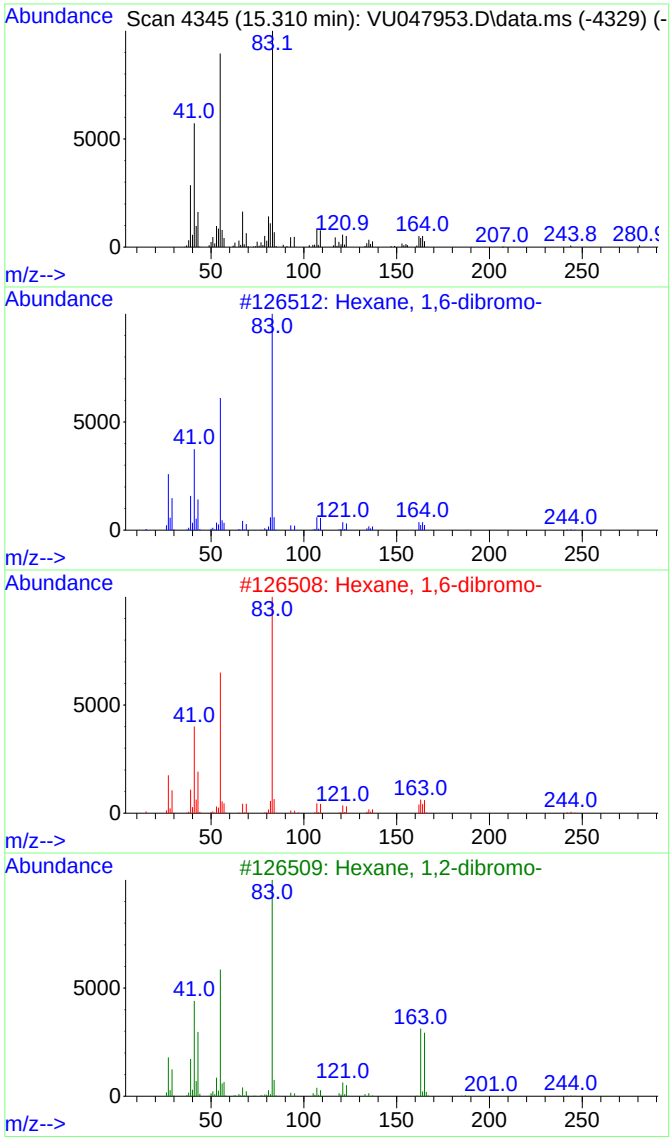
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM032422WMA.M
Quant Title : VOC Analysis

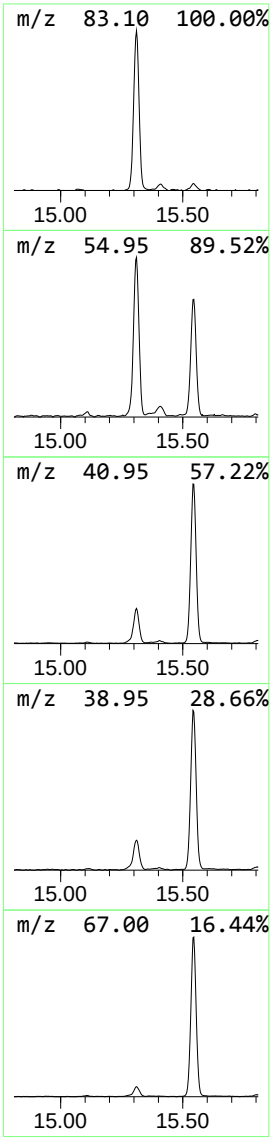
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Hexane, 1,6-dibromo- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.310	8.44 ug/L	231122	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 1,6-dibromo-	242	C6H12Br2	000629-03-8	97
2	Hexane, 1,6-dibromo-	242	C6H12Br2	000629-03-8	91
3	Hexane, 1,2-dibromo-	242	C6H12Br2	000624-20-4	72
4	Hexane, 1,5-dibromo-	242	C6H12Br2	000627-96-3	64
5	2-Butenoic acid, 2-methyl-, 2-me...	154	C9H14O2	061692-78-2	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041222\
Data File : VU047953.D
Acq On : 12 Apr 2022 14:17
Operator : SY/MD
Sample : N2293-08MEDL 5X
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ETNN8MEDL

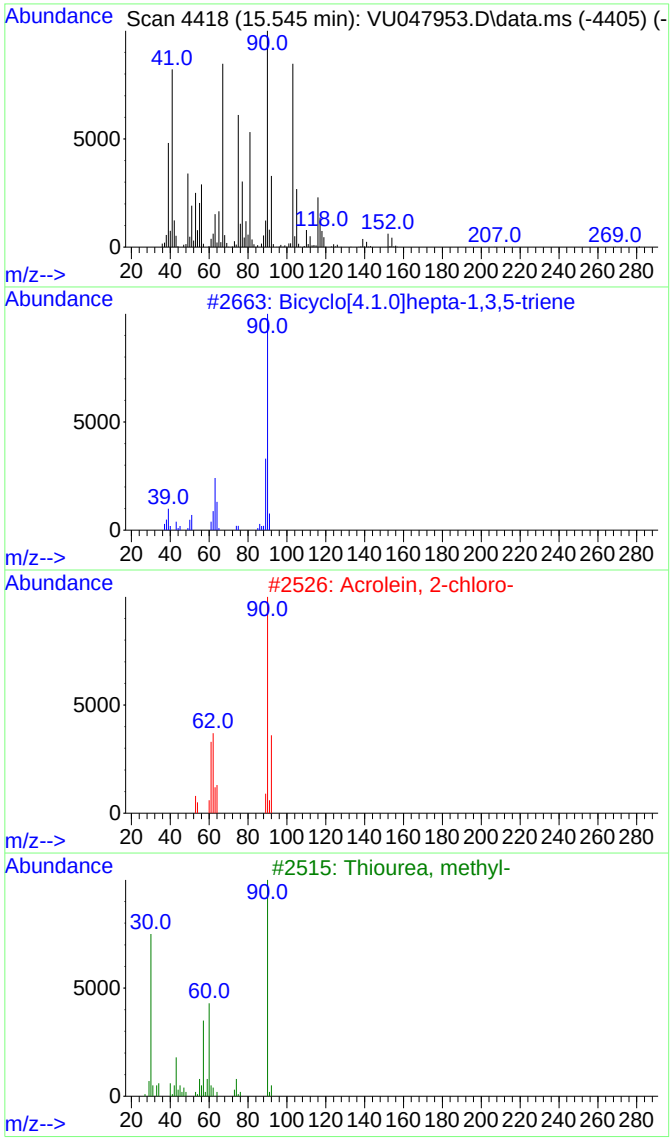
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM032422WMA.M
Quant Title : VOC Analysis

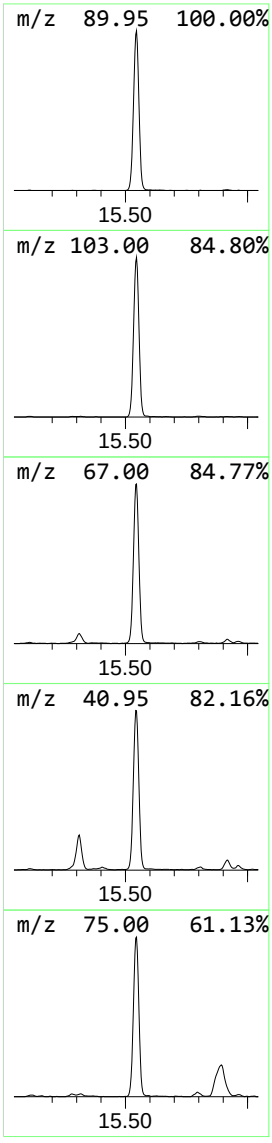
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.545	47.01 ug/L	1287850	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[4.1.0]hepta-1,3,5-triene	90	C7H6	004646-69-9	38
2	Acrolein, 2-chloro-	90	C3H3ClO	000683-51-2	37
3	Thiourea, methyl-	90	C2H6N2S	000598-52-7	35
4	Thiourea, methyl-	90	C2H6N2S	000598-52-7	35
5	1,3,5-Trichloro-4-methyl-2-penta...	202	C6H9Cl3O	1000140-03-8	25





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041222\
Data File : VU047953.D
Acq On : 12 Apr 2022 14:17
Operator : SY/MD
Sample : N2293-08MEDL 5X
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ETNN8MEDL

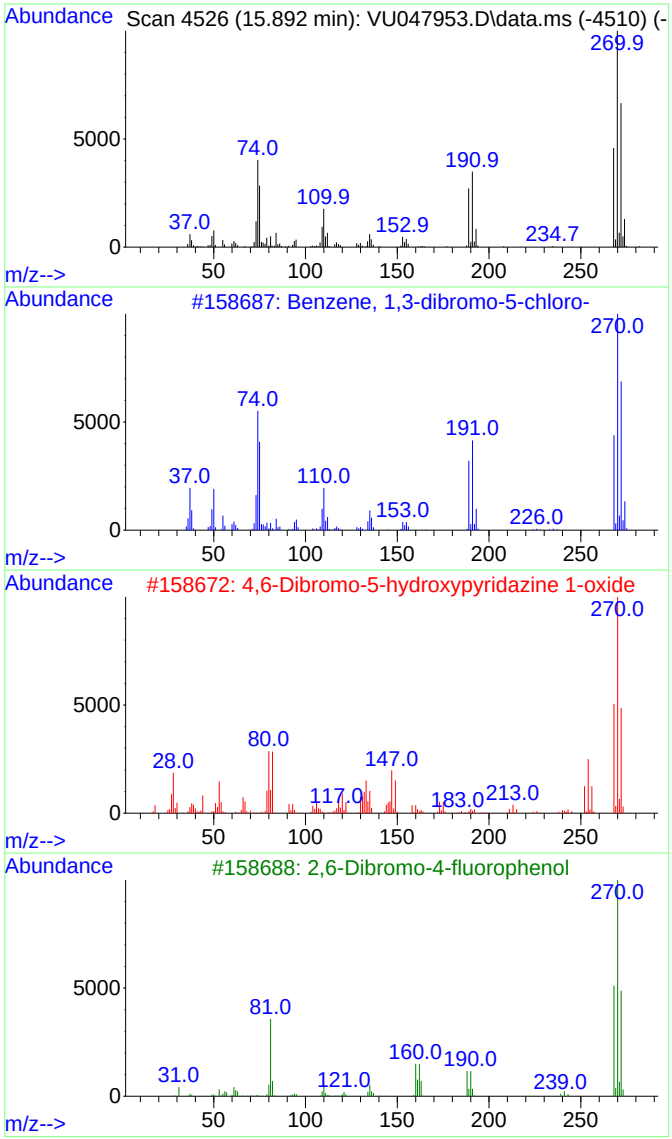
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM032422WMA.M
Quant Title : VOC Analysis

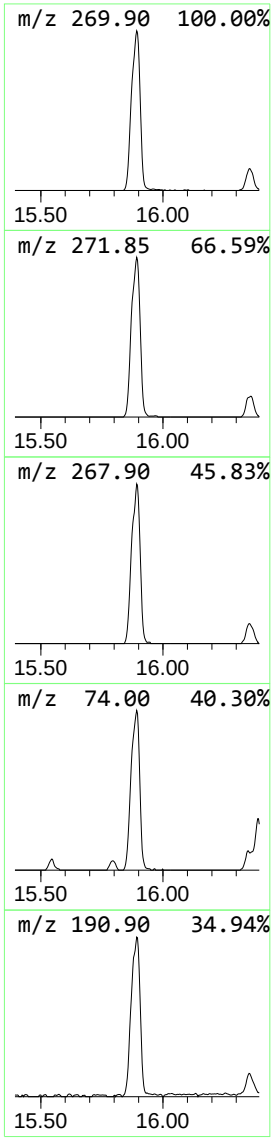
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1,3-dibromo-5-chloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.893	19.44 ug/L	532553	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dibromo-5-chloro-	268	C6H3Br2Cl	014862-52-3	94
2	4,6-Dibromo-5-hydroxypyridazine ...	268	C4H2Br2N2O2	019971-61-0	47
3	2,6-Dibromo-4-fluorophenol	268	C6H3Br2FO	000344-20-7	43
4	Pyrimidine, 4,6-dichloro-2-(meth...	270	C11H8Cl2N2S	064415-11-8	23
5	4-(4-Bromophenyl)-2-chloropyrimi...	268	C10H6BrClN2	932162-80-6	17





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041222\
Data File : VU047953.D
Acq On : 12 Apr 2022 14:17
Operator : SY/MD
Sample : N2293-08MEDL 5X
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ETNN8MEDL

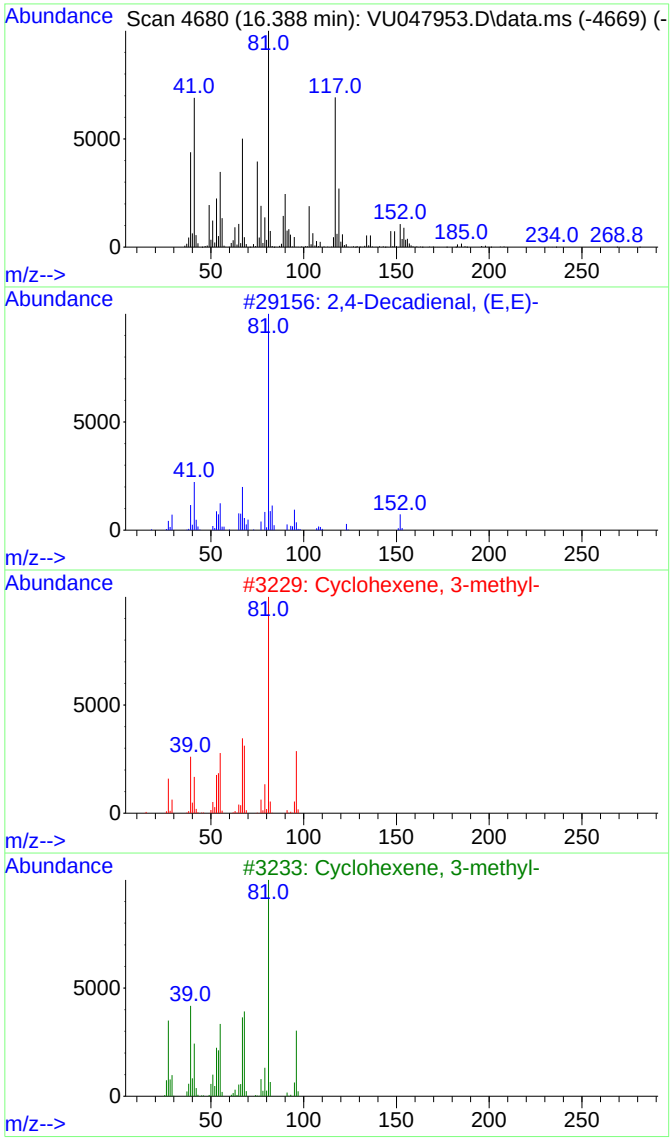
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM032422WMA.M
Quant Title : VOC Analysis

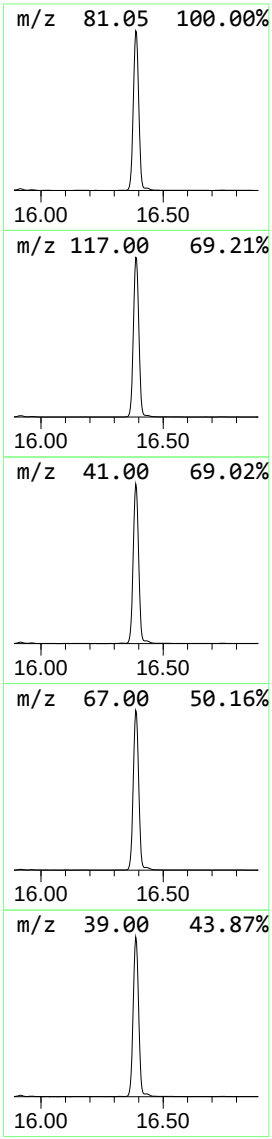
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 unknown-02 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.388	304.54 ug/L	8343280	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4-Decadienal, (E,E)-	152	C10H16O	025152-84-5	22
2	Cyclohexene, 3-methyl-	96	C7H12	000591-48-0	12
3	Cyclohexene, 3-methyl-	96	C7H12	000591-48-0	12
4	2-Hexyne, 4-methyl-	96	C7H12	020198-49-6	12
5	Cyclohexene, 1-methyl-	96	C7H12	000591-49-1	12





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041222\
Data File : VU047953.D
Acq On : 12 Apr 2022 14:17
Operator : SY/MD
Sample : N2293-08MEDL 5X
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ETNN8MEDL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM032422WMA.M
Quant Title : VOC Analysis

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

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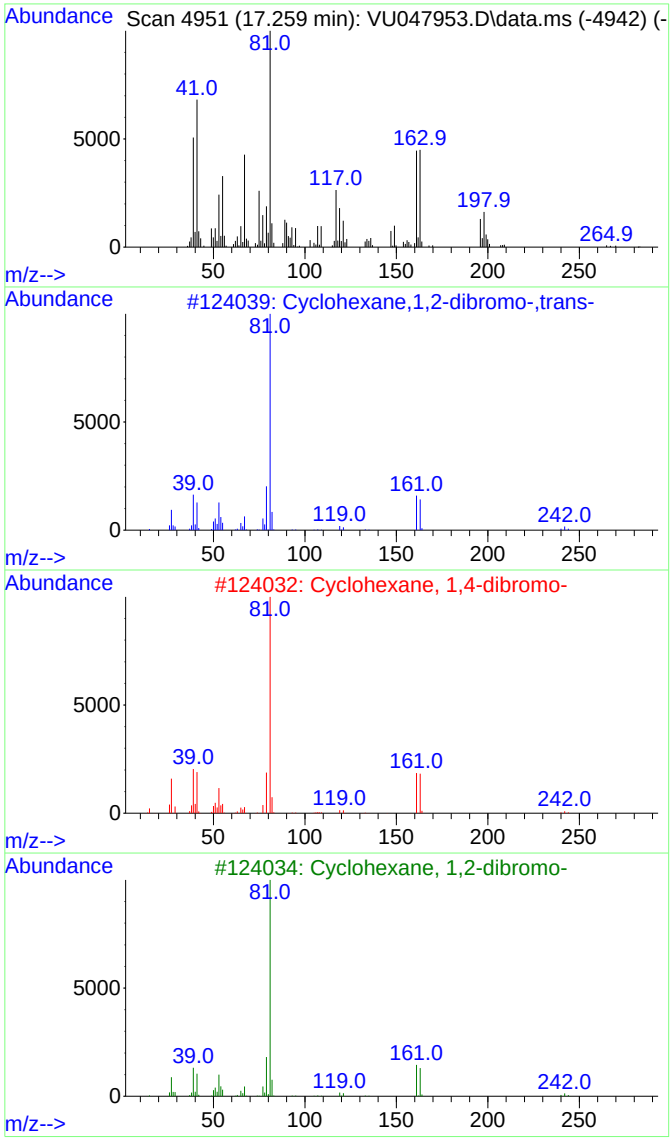
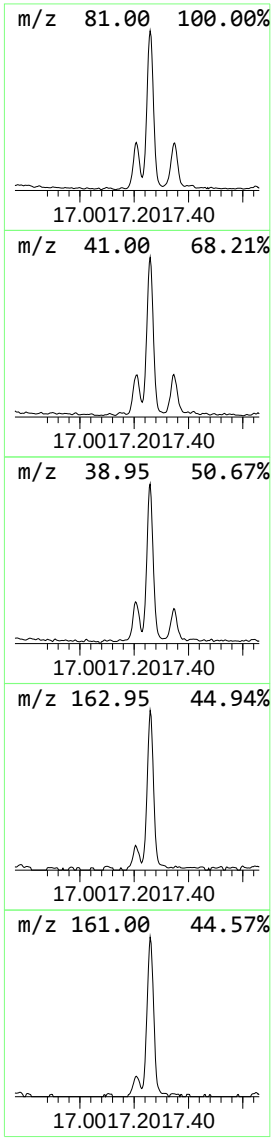
Peak Number 10 unknown-03 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.259	9.36 ug/L	256502	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane,1,2-dibromo-,trans-	240	C6H10Br2	007429-37-0	47
2	Cyclohexane, 1,4-dibromo-	240	C6H10Br2	035076-92-7	38
3	Cyclohexane, 1,2-dibromo-	240	C6H10Br2	005401-62-7	38
4	Methyl ethyl cyclopentene	110	C8H14	019780-56-4	35
5	Cyclopentaneethanol, 2-(hydroxym...	172	C10H20O2	000485-42-7	25





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041222\
Data File : VU047953.D
Acq On : 12 Apr 2022 14:17
Operator : SY/MD
Sample : N2293-08MEDL 5X
Misc : 6.28g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ETNN8MEDL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM032422WMA.M
Quant Title : VOC Analysis

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzene, 1-brom...	12.967	13.0	ug/L	357196	3	11.812	1369830	50.0
Benzene, 1-brom...	13.327	9.9	ug/L	270873	3	11.812	1369830	50.0
Hexane, 1,6-dic...	13.375	7.4	ug/L	203101	3	11.812	1369830	50.0
Hexane, 1-bromo...	14.359	7.7	ug/L	210575	3	11.812	1369830	50.0
Hexane, 1,6-dib...	15.310	8.4	ug/L	231122	3	11.812	1369830	50.0
unknown-01	15.545	47.0	ug/L	1287850	3	11.812	1369830	50.0
Benzene, 1,3-di...	15.893	19.4	ug/L	532553	3	11.812	1369830	50.0
unknown-02	16.388	304.5	ug/L	8343280	3	11.812	1369830	50.0
unknown-03	17.259	9.4	ug/L	256502	3	11.812	1369830	50.0