LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062822\
Data File : VU049463.D
Acq On : 28 Jun 2022 17:08
Operator : SY/MD
Sample : N3480-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
61522B

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_U\Method\82U062722W.M
Title : SW846 8260

Signal : TIC: VU049463.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.993	192	203	222	rVB	357237	486820	6.52%	0.937%
2	2.202	259	268	284	rVB2	56615	92615	1.24%	0.178%
3	2.295	284	297	323	rBV	154389	409474	5.48%	0.788%
4	2.626	378	400	429	rVB	328909	679687	9.10%	1.309%
5	2.768	432	444	466	rBV	44449	102358	1.37%	0.197%
6	3.041	513	529	533	rBV4	49892	97662	1.31%	0.188%
7	3.080	535	541	552	rVB2	79473	148483	1.99%	0.286%
8	3.363	612	629	651	rBV2	128274	286912	3.84%	0.552%
9	4.530	976	992	1012	rVB3	50184	127772	1.71%	0.246%
10	4.707	1036	1047	1075	rVB	37557	85558	1.15%	0.165%
11	5.288	1212	1228	1241	rBV	254112	538367	7.21%	1.036%
12	5.372	1241	1254	1273	rVB	468944	990771	13.27%	1.907%
13	5.703	1343	1357	1369	rBV	272989	554128	7.42%	1.067%
14	5.774	1369	1379	1393	rVB	138768	279138	3.74%	0.537%
15	5.899	1408	1418	1436	rVB2	39589	92750	1.24%	0.179%
16	6.250	1512	1527	1563	rVV	650302	1241303	16.63%	2.390%
17	6.764	1664	1687	1709	rBV3	56858	134868	1.81%	0.260%
18	7.793	1989	2007	2018	rBV2	39599	75897	1.02%	0.146%
19	7.896	2019	2039	2050	rVV	979114	1675986	22.45%	3.227%
20	7.970	2050	2062	2097	rVB	4385892	7465814	100.00%	14.373%
21	9.417	2499	2512	2530	rBV	882660	1433991	19.21%	2.761%
22	9.571	2547	2560	2578	rBV	617796	993554	13.31%	1.913%
23	9.690	2582	2597	2619	rBV	2540930	4213531	56.44%	8.112%
24	10.099	2710	2724	2751	rBV	1666993	2724205	36.49%	5.245%
25	10.484	2832	2844	2854	rBV	75037	118244	1.58%	0.228%
26	10.632	2878	2890	2906	rBV	678264	1076641	14.42%	2.073%
27	10.906	2963	2975	2987	rBV	151030	235573	3.16%	0.454%
28	10.996	2987	3003	3021	rVV	713418	1495695	20.03%	2.880%
29	11.086	3021	3031	3053	rVB	410954	660153	8.84%	1.271%
30	11.291	3083	3095	3114	rVB	462823	727501	9.74%	1.401%
31	11.465	3137	3149	3163	rBV	1384119	2150214	28.80%	4.140%
32	11.642	3197	3204	3216	rVB	52234	81363	1.09%	0.157%
33	11.742	3219	3235	3244	rBV	79912	125996	1.69%	0.243%
34	11.809	3244	3256	3270	rVV	863556	1418601	19.00%	2.731%
35	11.893	3270	3282	3297	rVV	689823	1078225	14.44%	2.076%
36	12.060	3324	3334	3337	rBV	114998	168847	2.26%	0.325%

LSC Area Percent Report

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Sample : N3480-03
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ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
61522B

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

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Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Title : SW846 8260

37	12.095	3337	3345	3351	rVV2	325592	476119	6.38%	0.917%
38	12.189	3363	3374	3390	rVB3	298828	573185	7.68%	1.103%
39	12.304	3399	3410	3423	rBV5	49346	109395	1.47%	0.211%
40	12.375	3423	3432	3446	rVB	144974	229331	3.07%	0.442%
41	12.465	3447	3460	3465	rBV	215370	350975	4.70%	0.676%
42	12.494	3465	3469	3480	rVB	207351	281264	3.77%	0.541%
43	12.571	3481	3493	3500	rBV	299837	456764	6.12%	0.879%
44	12.610	3500	3505	3516	rVB	77457	113471	1.52%	0.218%
45	12.680	3516	3527	3537	rBV	219731	439617	5.89%	0.846%
46	12.873	3577	3587	3599	rVB	149251	246819	3.31%	0.475%
47	12.970	3599	3617	3625	rBV	228454	377247	5.05%	0.726%
48	13.024	3625	3634	3650	rVB	425975	658630	8.82%	1.268%
49	13.108	3650	3660	3664	rBV3	52297	85066	1.14%	0.164%
50	13.291	3709	3717	3727	rVB2	341187	534160	7.15%	1.028%
51	13.352	3729	3736	3744	rVB3	54079	75681	1.01%	0.146%
52	13.404	3744	3752	3756	rBV2	77288	118663	1.59%	0.228%
53	13.452	3757	3767	3780	rVB3	976967	1702493	22.80%	3.278%
54	13.623	3809	3820	3843	rVB	810310	1314487	17.61%	2.531%
55	13.732	3844	3854	3862	rBV3	82001	142607	1.91%	0.275%
56	13.783	3862	3870	3878	rVV	155617	269106	3.60%	0.518%
57	13.835	3878	3886	3901	rVV5	213008	504550	6.76%	0.971%
58	13.909	3904	3909	3913	rVV2	95622	147321	1.97%	0.284%
59	13.941	3913	3919	3935	rVB2	199267	394641	5.29%	0.760%
60	14.086	3951	3964	3971	rBV	153603	264825	3.55%	0.510%
61	14.188	3985	3996	4010	rVB	271882	461240	6.18%	0.888%
62	14.285	4010	4026	4037	rBV2	319712	579560	7.76%	1.116%
63	14.346	4038	4045	4054	rVB2	103947	158946	2.13%	0.306%
64	14.484	4080	4088	4111	rVB	238519	392188	5.25%	0.755%
65	14.684	4133	4150	4169	rBV	816235	1646468	22.05%	3.170%
66	14.809	4180	4189	4198	rBV2	99169	159420	2.14%	0.307%
67	14.877	4200	4210	4221	rVB3	234569	394532	5.28%	0.760%
68	15.015	4242	4253	4267	rBV2	622906	1023580	13.71%	1.971%
69	15.198	4300	4310	4313	rBV2	245714	362247	4.85%	0.697%
70	15.269	4326	4332	4346	rVB2	163127	282263	3.78%	0.543%
71	15.346	4347	4356	4364	rBV3	58177	94175	1.26%	0.181%
72	15.407	4365	4375	4385	rVV3	239032	427045	5.72%	0.822%
73	15.465	4385	4393	4399	rVV3	107540	195557	2.62%	0.376%
74	15.500	4400	4404	4415	rVB2	82365	122460	1.64%	0.236%
75	15.574	4418	4427	4442	rVV3	98835	225224	3.02%	0.434%
76	15.764	4474	4486	4496	rBV4	52002	90719	1.22%	0.175%

LSC Area Percent Report

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ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
61522B

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
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Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Title : SW846 8260

77	15.925	4526	4536	4555	rBV3	223432	475056	6.36%	0.915%
78	16.147	4597	4605	4612	rBV3	64116	102612	1.37%	0.198%
79	16.262	4632	4641	4648	rBV3	114710	194703	2.61%	0.375%
80	16.407	4677	4686	4693	rBV	154982	256617	3.44%	0.494%
81	16.446	4693	4698	4714	rVB2	98606	162956	2.18%	0.314%

Sum of corrected areas: 51942682

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062822\
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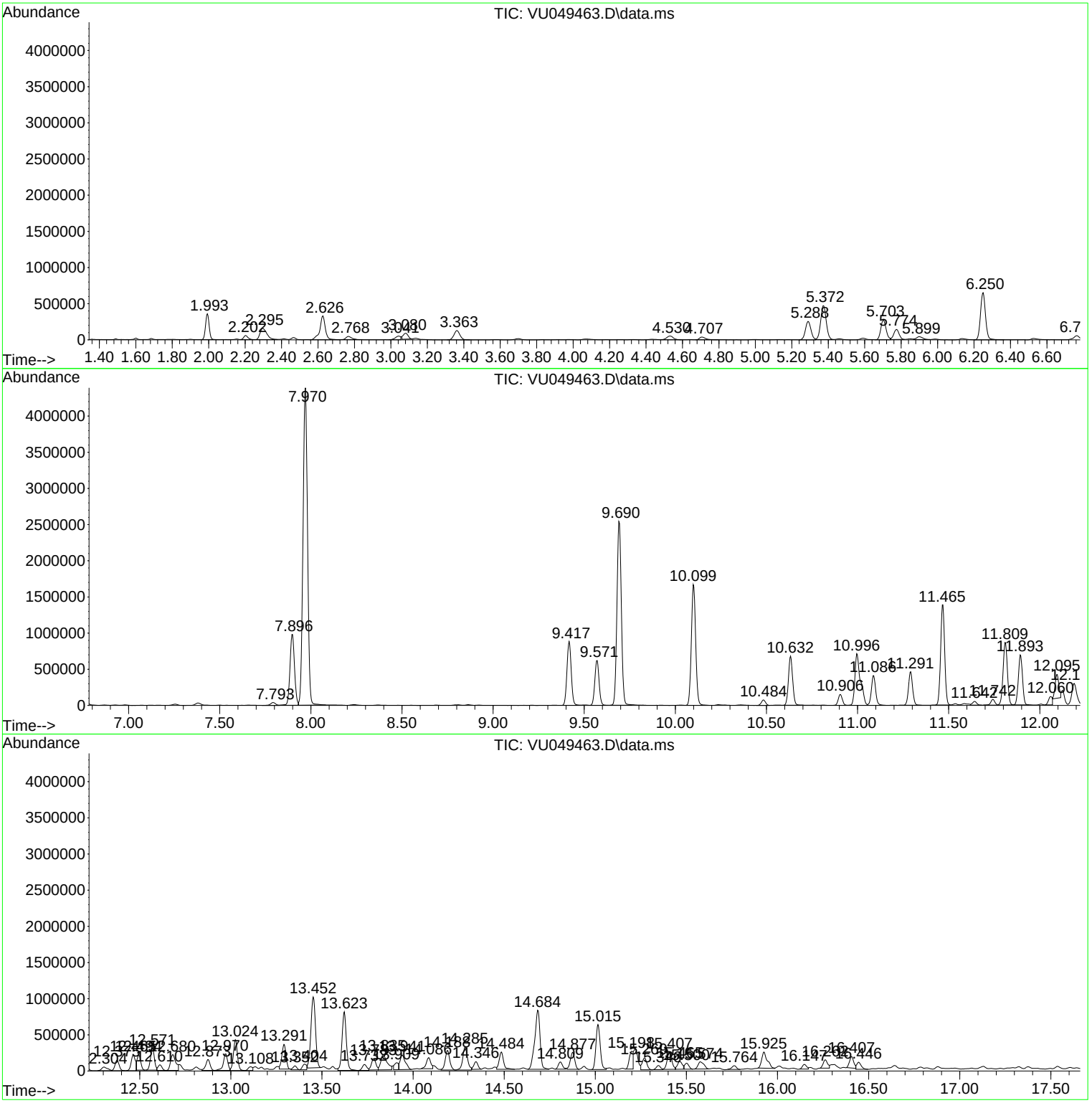
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
61522B

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062822\
Data File : VU049463.D
Acq On : 28 Jun 2022 17:08
Operator : SY/MD
Sample : N3480-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

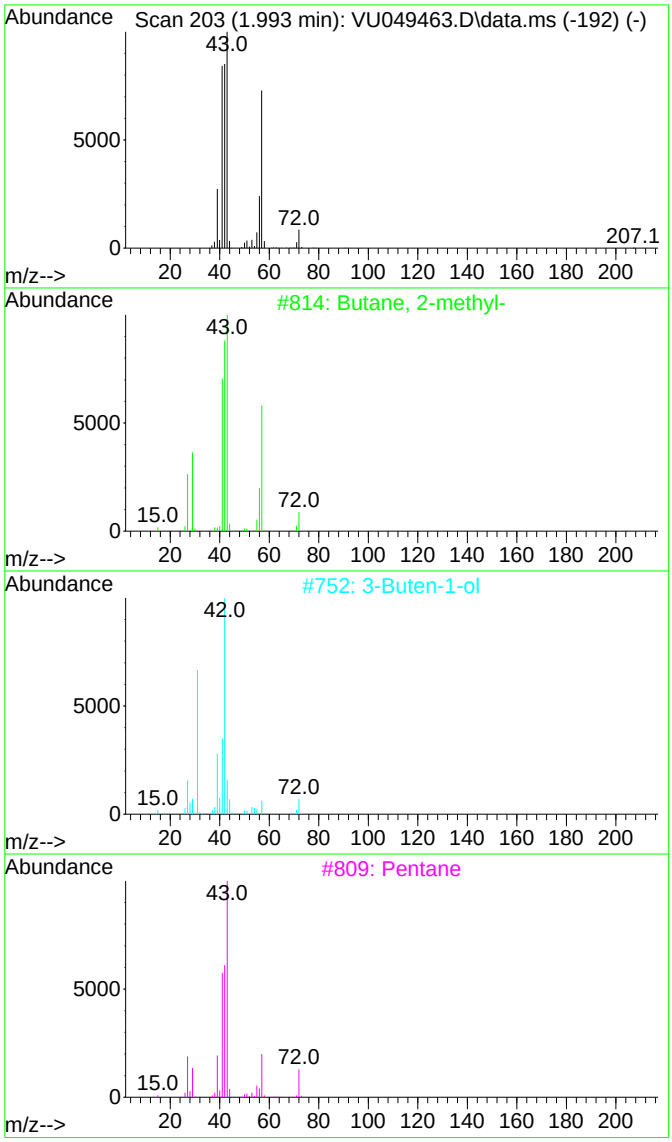
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

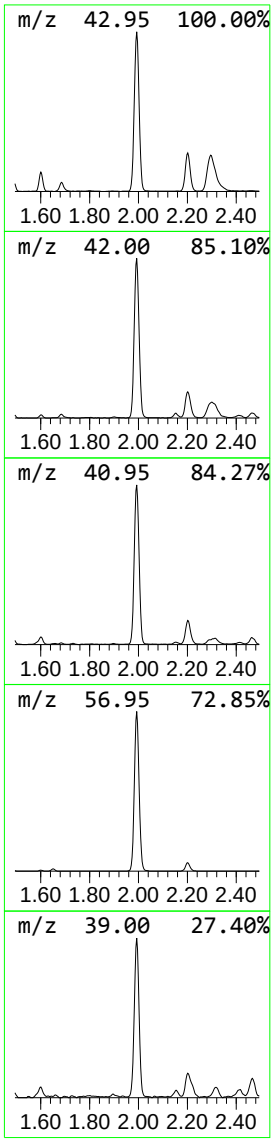
****

Peak Number 1 Butane, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.993	24.57 ug/l	486820	Pentafluorobenzene	5.372

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	87
2	3-Buten-1-ol	72	C4H8O	000627-27-0	47
3	Pentane	72	C5H12	000109-66-0	36
4	1-Butene	56	C4H8	000106-98-9	10
5	1-Propene, 2-methyl-	56	C4H8	000115-11-7	9





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062822\
Data File : VU049463.D
Acq On : 28 Jun 2022 17:08
Operator : SY/MD
Sample : N3480-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
61522B

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

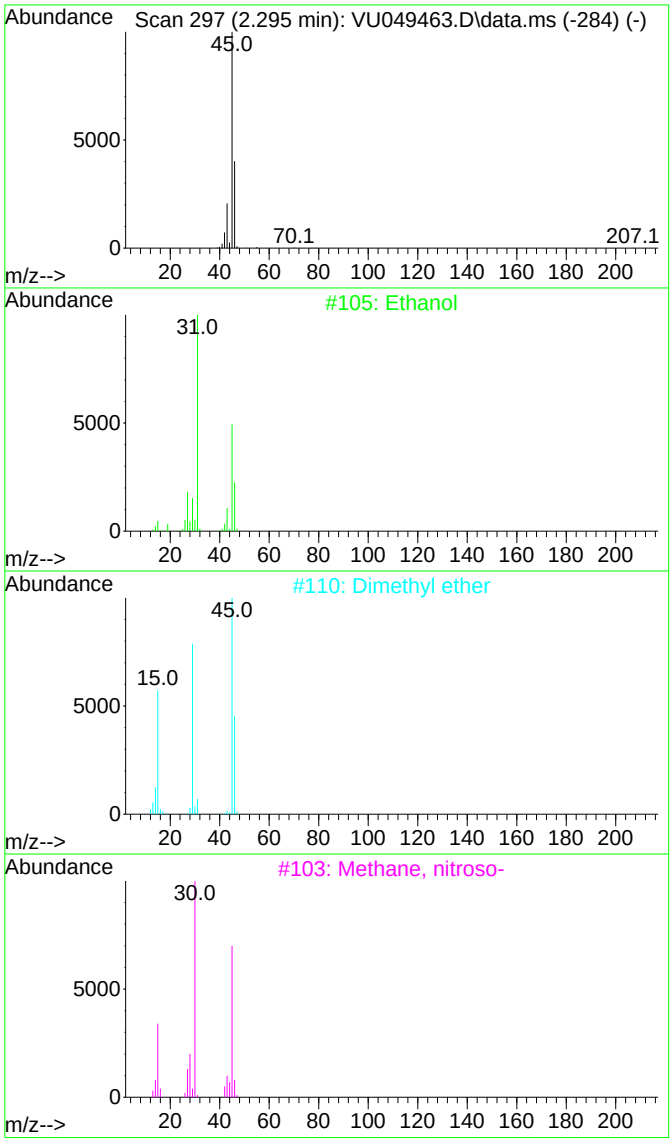
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

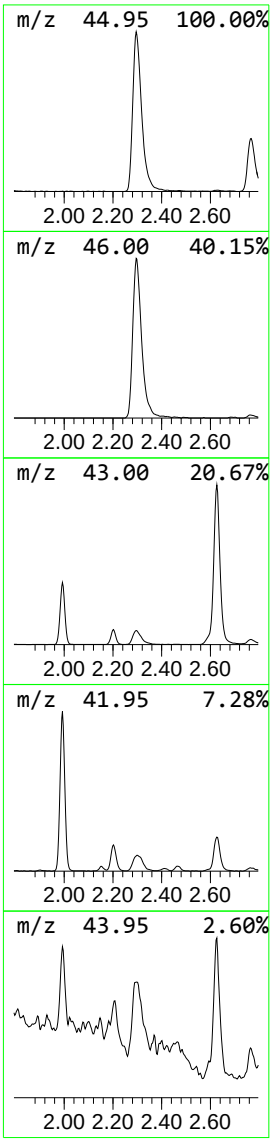
****

Peak Number 2 Ethanol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.295	20.66 ug/l	409474	Pentafluorobenzene	5.372

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethanol	46	C2H6O	000064-17-5	64
2	Dimethyl ether	46	C2H6O	000115-10-6	9
3	Methane, nitroso-	45	CH3NO	000865-40-7	4
4	Formic acid	46	CH2O2	000064-18-6	4
5	Oxalic acid	90	C2H2O4	000144-62-7	4





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062822\
Data File : VU049463.D
Acq On : 28 Jun 2022 17:08
Operator : SY/MD
Sample : N3480-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
61522B

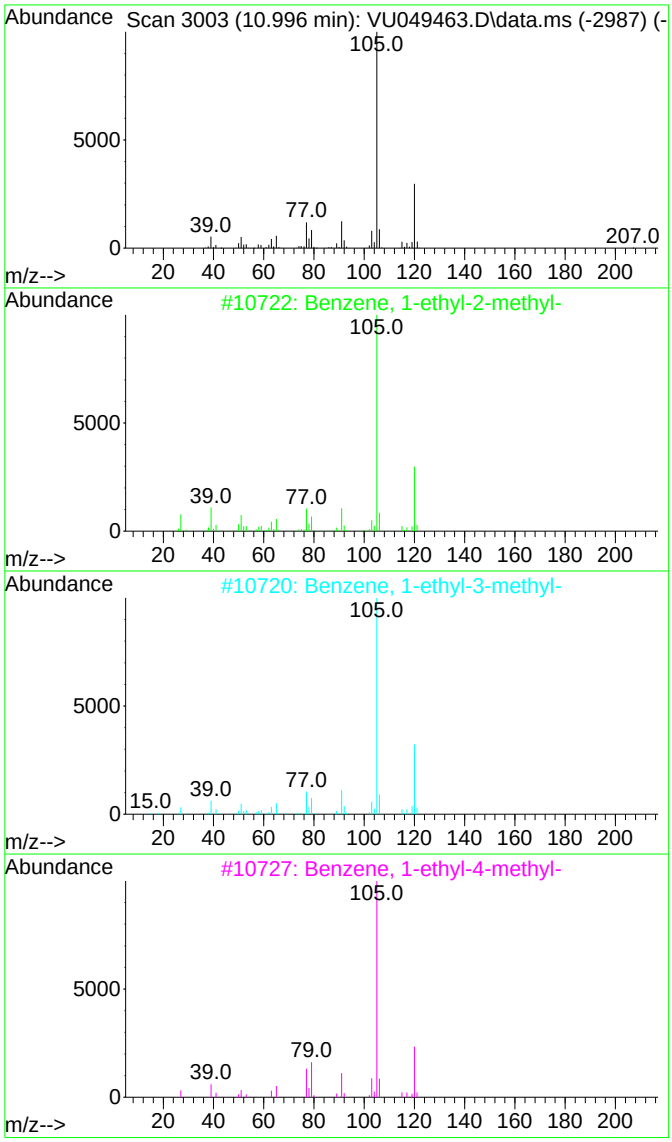
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

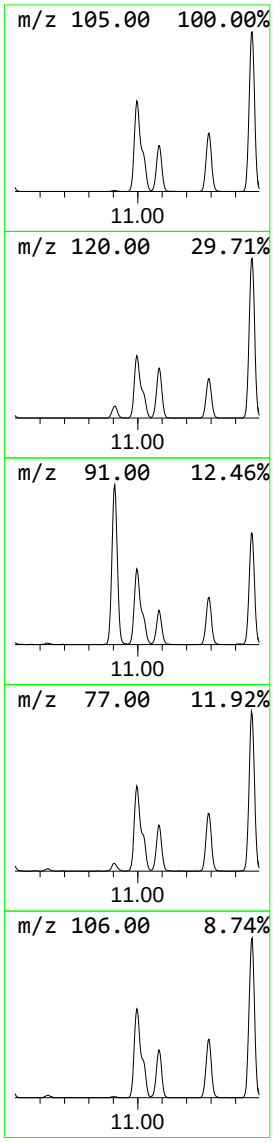
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.996	52.72 ug/l	1495700	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Mesitylene	120	C9H12	000108-67-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062822\
Data File : VU049463.D
Acq On : 28 Jun 2022 17:08
Operator : SY/MD
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Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
61522B

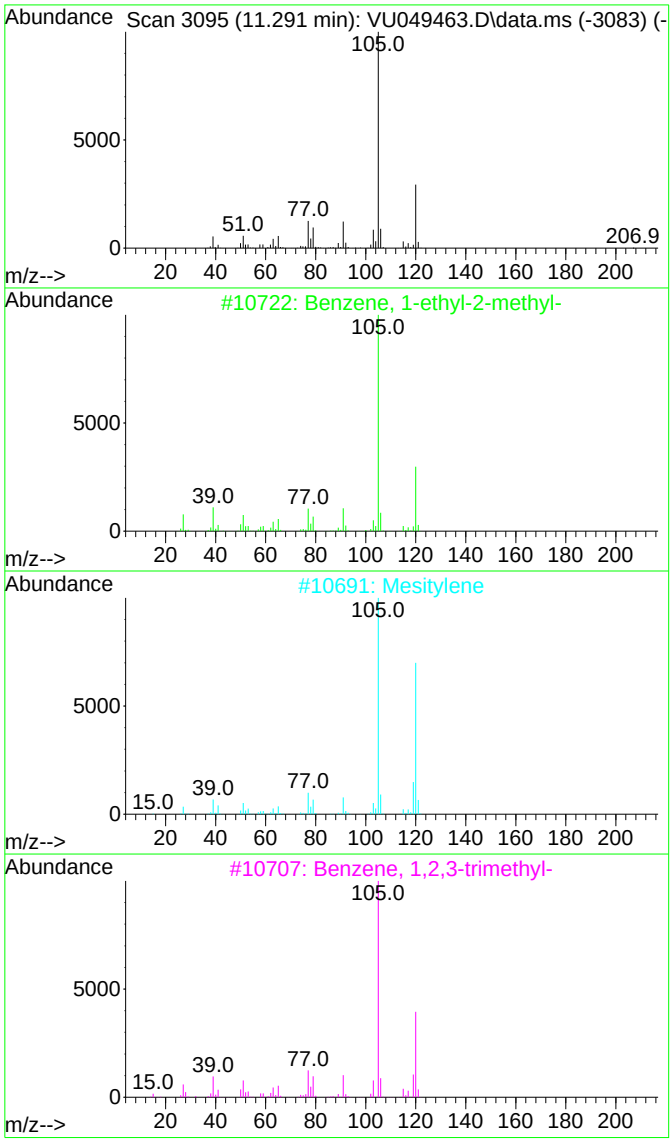
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

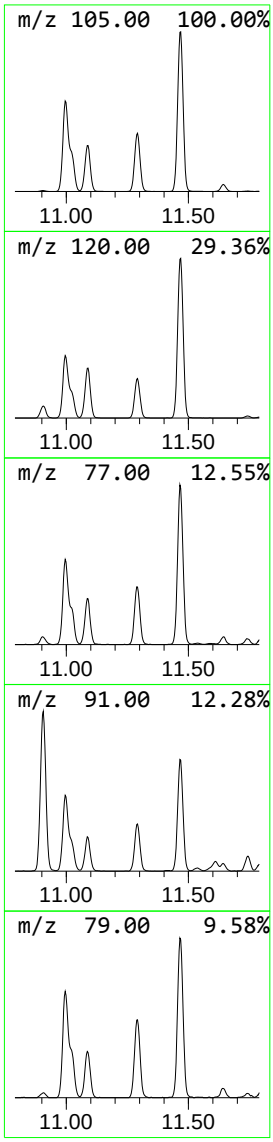
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 1,2,3-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.291	25.64 ug/l	727501	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2	Mesitylene	120	C9H12	000108-67-8	91
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





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Acq On : 28 Jun 2022 17:08
Operator : SY/MD
Sample : N3480-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
61522B

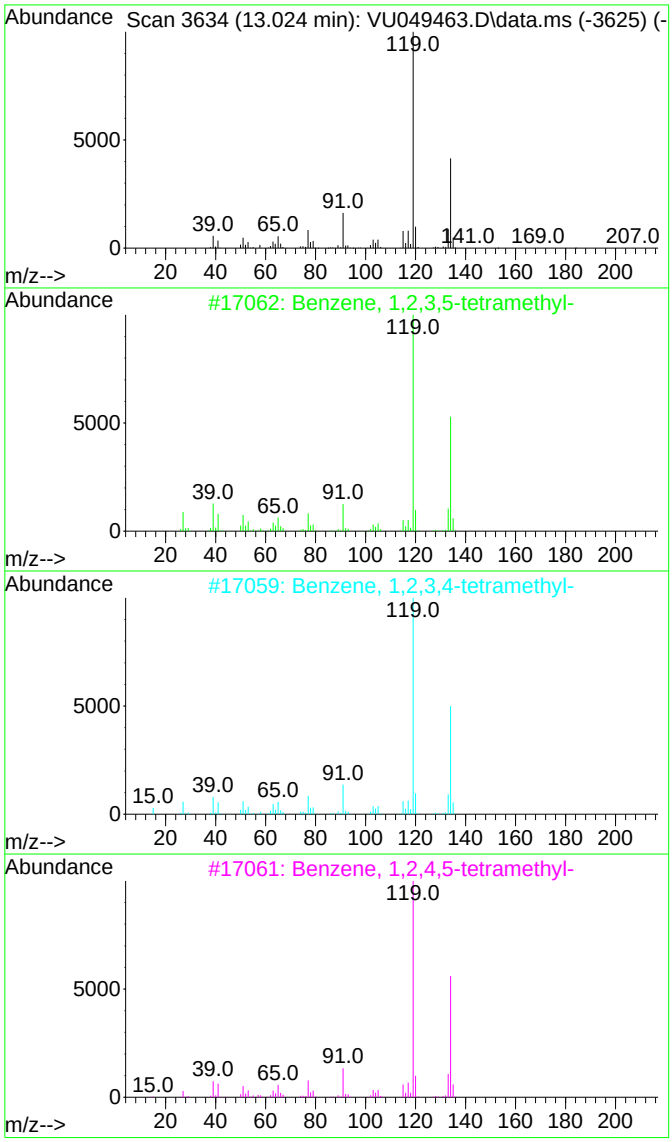
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

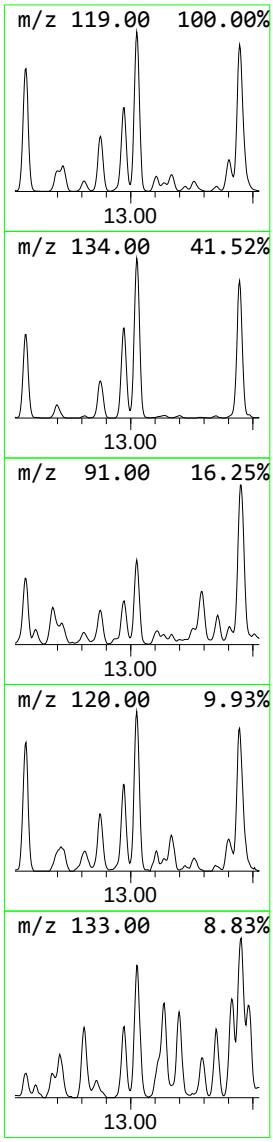
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.025	23.21 ug/l	658630	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062822\
Data File : VU049463.D
Acq On : 28 Jun 2022 17:08
Operator : SY/MD
Sample : N3480-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
61522B

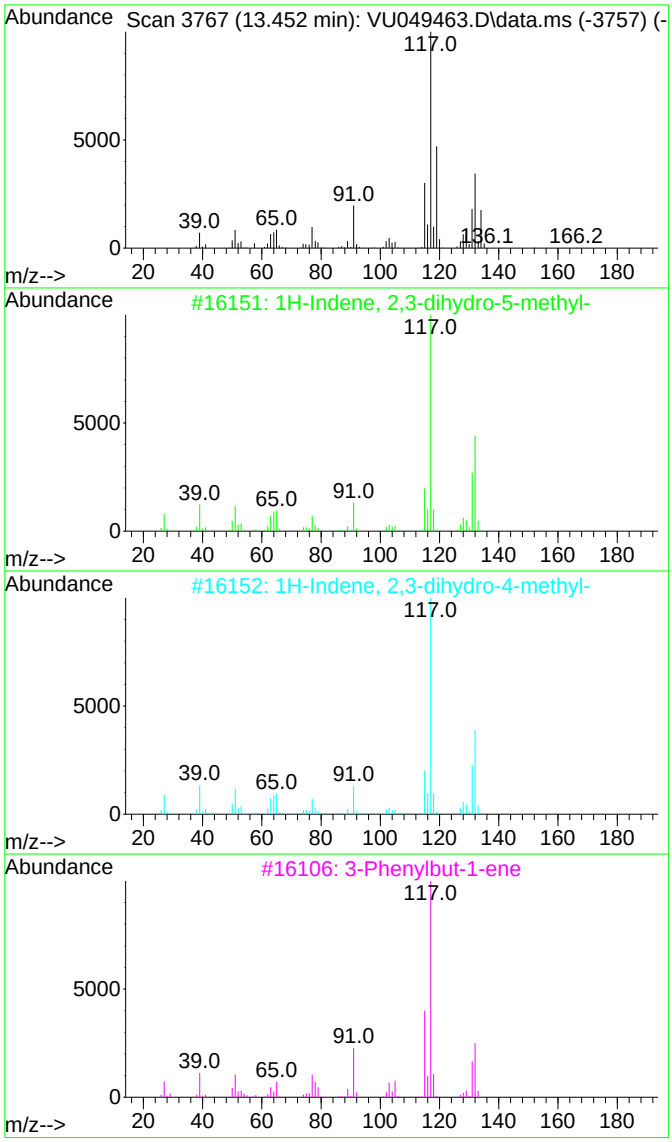
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

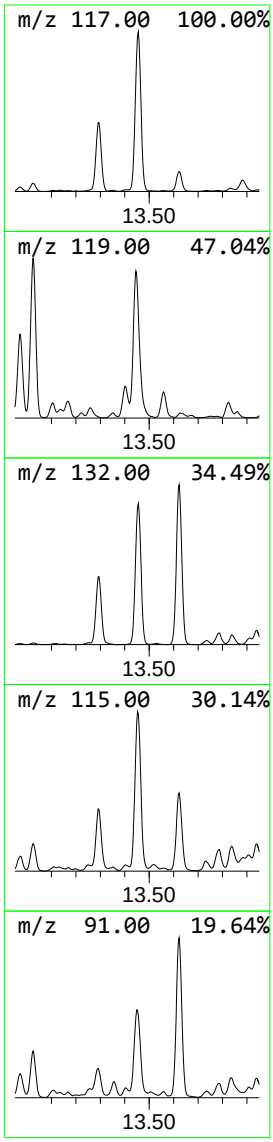
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.452	60.01 ug/l	1702490	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
2	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	92
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
4	Indan, 1-methyl-	132	C10H12	000767-58-8	70
5	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	70





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062822\
Data File : VU049463.D
Acq On : 28 Jun 2022 17:08
Operator : SY/MD
Sample : N3480-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
61522B

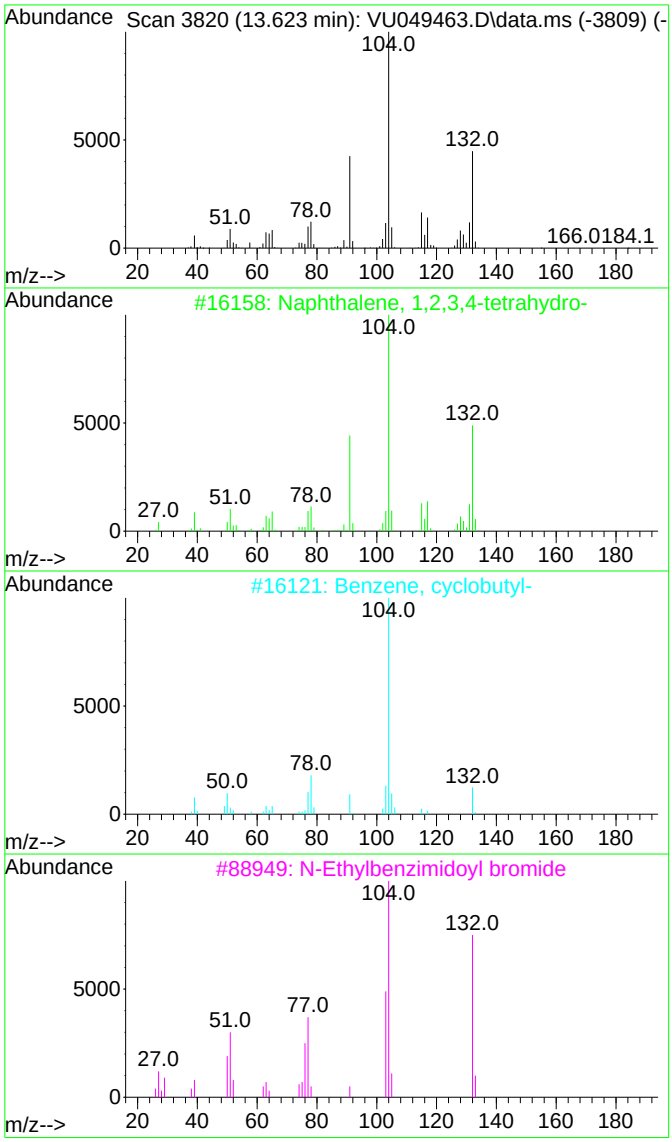
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

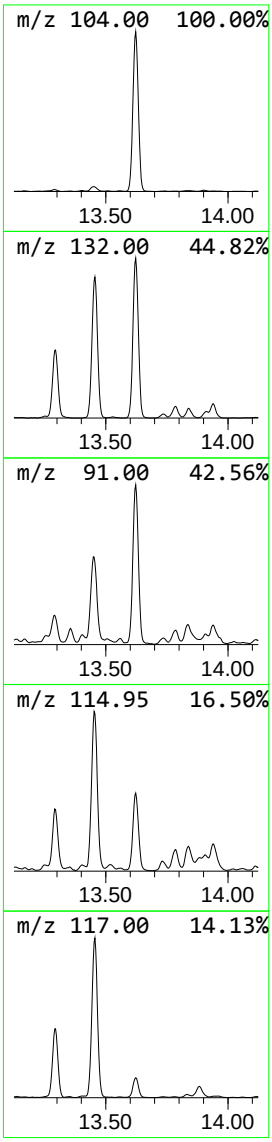
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.623	46.33 ug/l	1314490	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	95
2	Benzene, cyclobutyl-	132	C10H12	004392-30-7	42
3	N-Ethylbenzimidoyl bromide	211	C9H10BrN	041182-87-0	40
4	3(2H)-Furanone, dihydro-2,2-dime...	190	C12H14O2	063678-00-2	38
5	Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	020071-09-4	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062822\
Data File : VU049463.D
Acq On : 28 Jun 2022 17:08
Operator : SY/MD
Sample : N3480-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
61522B

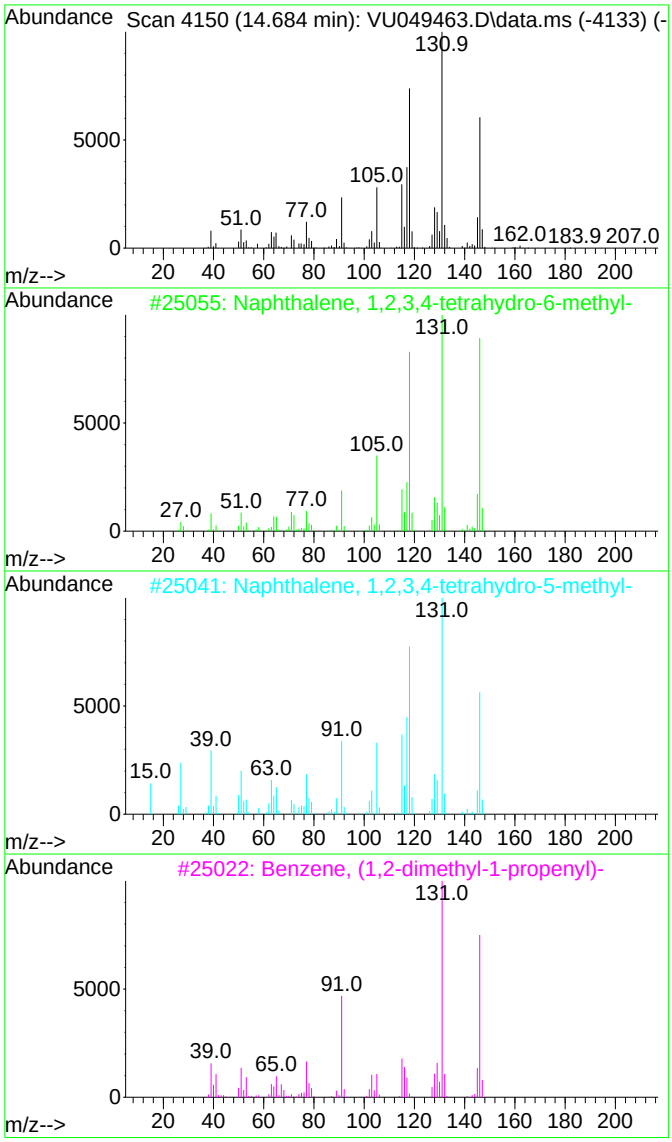
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

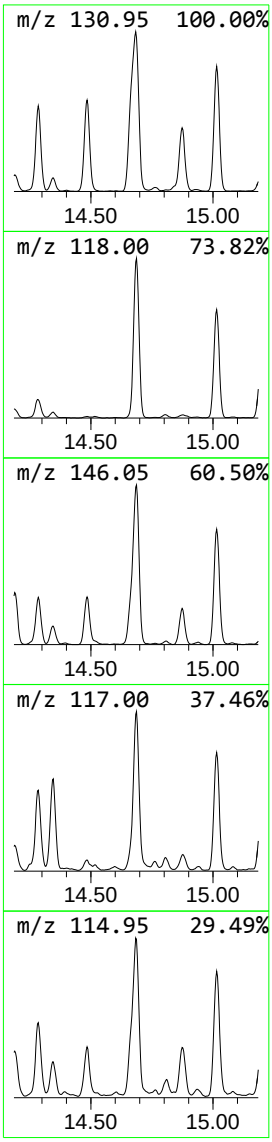
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.684	58.03 ug/l	1646470	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	97
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	96
3	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	90
4	Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	78
5	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	60





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062822\
Data File : VU049463.D
Acq On : 28 Jun 2022 17:08
Operator : SY/MD
Sample : N3480-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
61522B

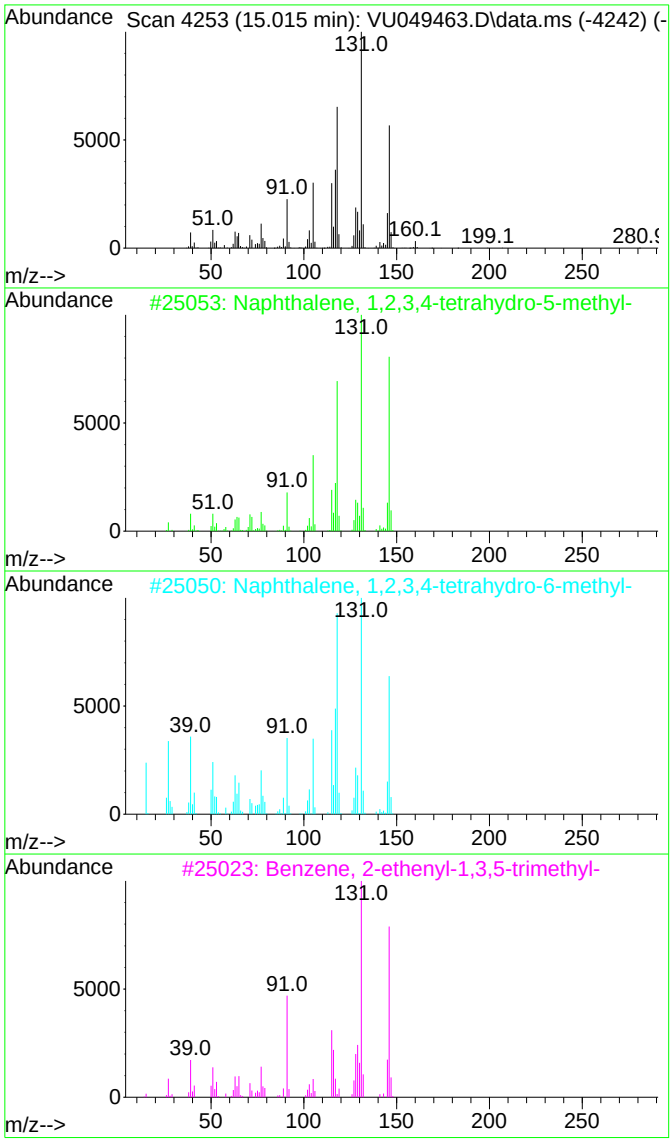
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

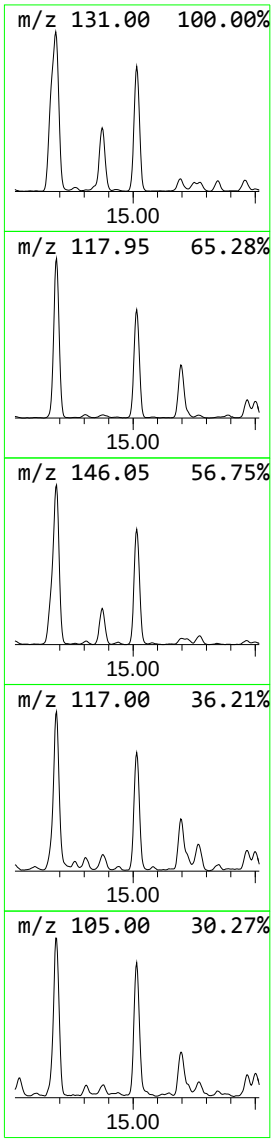
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.015	36.08 ug/l	1023580	1,4-Dichlorobenzene-d4	11.812

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	96
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	96
3	Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	86
4	1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	74
5	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	70





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062822\
Data File : VU049463.D
Acq On : 28 Jun 2022 17:08
Operator : SY/MD
Sample : N3480-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
61522B

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-methyl-	1.993	24.6	ug/l	486820	1	5.372	990771	50.0
Ethanol	2.295	20.7	ug/l	409474	1	5.372	990771	50.0
Benzene, 1-ethy...	10.996	52.7	ug/l	1495700	4	11.812	1418600	50.0
Benzene, 1,2,3-...	11.291	25.6	ug/l	727501	4	11.812	1418600	50.0
Benzene, 1,2,3,...	13.025	23.2	ug/l	658630	4	11.812	1418600	50.0
1H-Indene, 2,3-...	13.452	60.0	ug/l	1702490	4	11.812	1418600	50.0
Naphthalene, 1,...	13.623	46.3	ug/l	1314490	4	11.812	1418600	50.0
Naphthalene, 1,...	14.684	58.0	ug/l	1646470	4	11.812	1418600	50.0
Naphthalene, 1,...	15.015	36.1	ug/l	1023580	4	11.812	1418600	50.0