LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060053.D
Acq On : 19 Jul 2024 11:50
Operator : MD/SY
Sample : P3217-03DL 100X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZC5DL

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.1 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\SFAMUTR071824WMA.M
Title : TRACE VOA SFAM1.0

Signal : TIC: VU060053.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.596	85	95	109	rBV2	52216	63254	6.63%	0.908%
2	1.911	184	193	212	rVB	45313	62156	6.51%	0.893%
3	2.563	384	396	413	rBV	75711	124454	13.04%	1.787%
4	2.679	424	432	444	rVB3	5530	10270	1.08%	0.147%
5	4.657	1026	1047	1085	rBV2	356654	954696	100.00%	13.710%
6	5.055	1156	1171	1194	rBV2	82527	185441	19.42%	2.663%
7	5.721	1356	1378	1403	rBV3	160246	428831	44.92%	6.158%
8	6.242	1527	1540	1561	rBV	157344	308484	32.31%	4.430%
9	6.541	1621	1633	1652	rBV2	15417	31422	3.29%	0.451%
10	6.685	1665	1678	1696	rBV	103497	198559	20.80%	2.852%
11	7.560	1939	1950	1970	rBV2	44759	81496	8.54%	1.170%
12	7.891	2040	2053	2069	rBV	186232	326327	34.18%	4.686%
13	7.968	2069	2077	2090	rVB	13565	25239	2.64%	0.362%
14	8.177	2131	2142	2160	rBV2	25455	45965	4.81%	0.660%
15	8.547	2247	2257	2271	rVB	34138	58174	6.09%	0.835%
16	8.627	2271	2282	2319	rBV	206867	415949	43.57%	5.973%
17	9.412	2514	2526	2548	rBV	234971	404323	42.35%	5.806%
18	9.566	2565	2574	2587	rVB	25694	40109	4.20%	0.576%
19	9.688	2600	2612	2638	rBV	66136	116008	12.15%	1.666%
20	10.094	2724	2738	2759	rBV	86073	144908	15.18%	2.081%
21	10.750	2931	2942	2954	rBV	88302	142468	14.92%	2.046%
22	10.894	2970	2987	3002	rBV4	10081	22238	2.33%	0.319%
23	10.994	3008	3018	3023	rBV2	34984	56757	5.95%	0.815%
24	11.084	3036	3046	3064	rVB	46992	74193	7.77%	1.065%
25	11.286	3098	3109	3122	rBV	46629	73336	7.68%	1.053%
26	11.463	3153	3164	3182	rBV	75837	121240	12.70%	1.741%
27	11.631	3202	3216	3228	rVB8	7192	13205	1.38%	0.190%
28	11.740	3243	3250	3257	rBV3	9934	13126	1.37%	0.189%
29	11.807	3257	3271	3287	rVV	250379	469565	49.18%	6.743%
30	11.888	3287	3296	3309	rVB	119413	189499	19.85%	2.721%
31	12.090	3348	3359	3366	rBV3	44588	76679	8.03%	1.101%
32	12.190	3377	3390	3414	rVB2	273091	618460	64.78%	8.882%
33	12.296	3415	3423	3437	rVB7	5769	10683	1.12%	0.153%
34	12.373	3437	3447	3462	rVB2	17371	26444	2.77%	0.380%
35	12.460	3464	3474	3479	rBV2	20469	31979	3.35%	0.459%
36	12.489	3479	3483	3497	rVB	20750	29022	3.04%	0.417%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060053.D
Acq On : 19 Jul 2024 11:50
Operator : MD/SY
Sample : P3217-03DL 100X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZC5DL

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.1 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\SFAMUTR071824WMA.M
Title : TRACE VOA SFAM1.0

37	12.566	3497	3507	3514	rBV2	39780	60276	6.31%	0.866%
38	12.605	3514	3519	3529	rVB2	10856	15866	1.66%	0.228%
39	12.676	3531	3541	3560	rBV2	29215	65327	6.84%	0.938%
40	12.868	3590	3601	3615	rVV3	19098	32105	3.36%	0.461%
41	12.968	3615	3632	3640	rVV3	25991	48775	5.11%	0.700%
42	13.020	3640	3648	3662	rVB	42908	65144	6.82%	0.936%
43	13.290	3724	3732	3740	rVB	16667	24401	2.56%	0.350%
44	13.447	3770	3781	3794	rVV4	69473	123278	12.91%	1.770%
45	13.508	3794	3800	3811	rVB5	10836	16915	1.77%	0.243%
46	13.618	3824	3834	3847	rVB4	28914	48189	5.05%	0.692%
47	13.730	3860	3869	3877	rBV6	7595	11588	1.21%	0.166%
48	13.836	3892	3902	3914	rVV7	11468	21775	2.28%	0.313%
49	13.933	3926	3932	3948	rVB4	8354	17436	1.83%	0.250%
50	14.084	3964	3979	3997	rBV	38753	70592	7.39%	1.014%
51	14.283	4030	4041	4050	rBV7	8468	14700	1.54%	0.211%
52	14.479	4093	4102	4116	rVB3	13591	21137	2.21%	0.304%
53	14.573	4120	4131	4147	rBV2	10295	17369	1.82%	0.249%
54	14.666	4147	4160	4174	rBV4	17955	43682	4.58%	0.627%
55	14.868	4215	4223	4233	rVB2	14322	21841	2.29%	0.314%
56	15.013	4255	4268	4275	rBV7	7193	12210	1.28%	0.175%
57	15.219	4317	4332	4346	rBV	76603	132375	13.87%	1.901%
58	15.405	4377	4390	4401	rBV	50220	83373	8.73%	1.197%

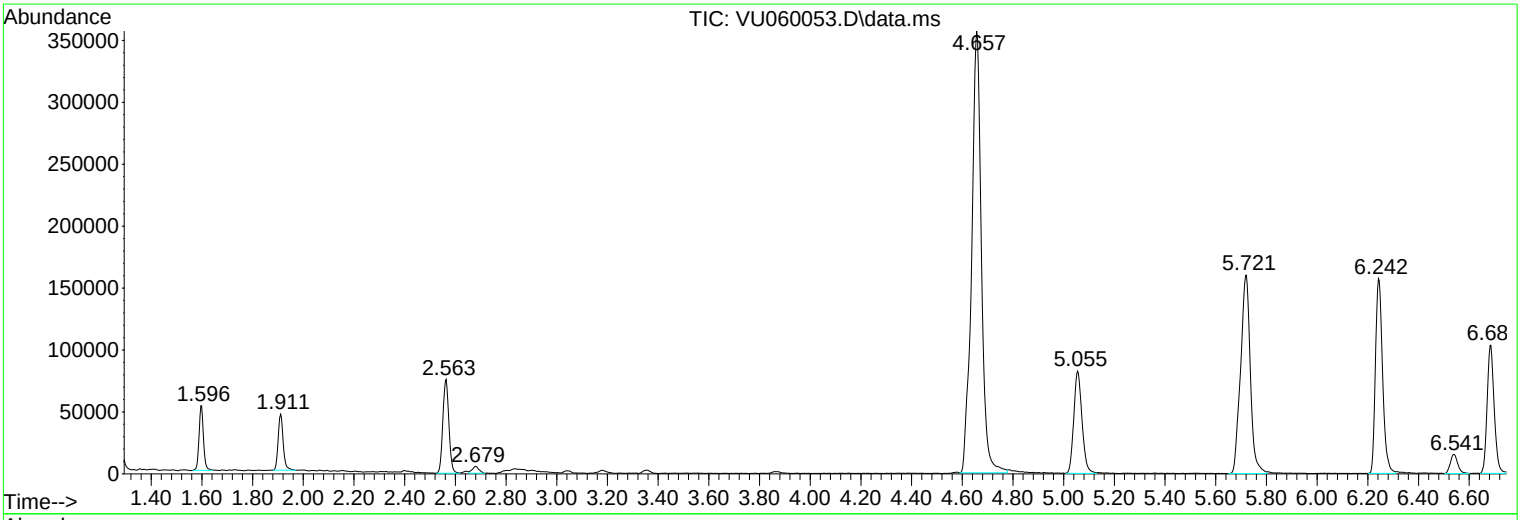
Sum of corrected areas: 6963313

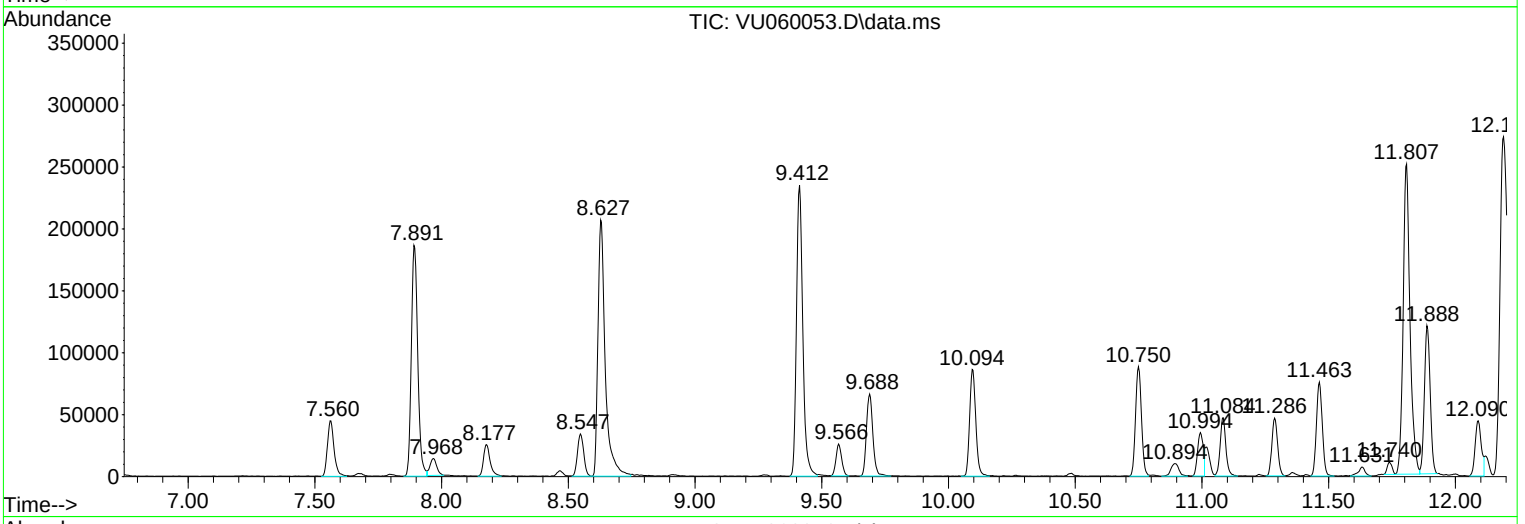
Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060053.D
Acq On : 19 Jul 2024 11:50
Operator : MD/SY
Sample : P3217-03DL 100X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 48 Sample Multiplier: 1

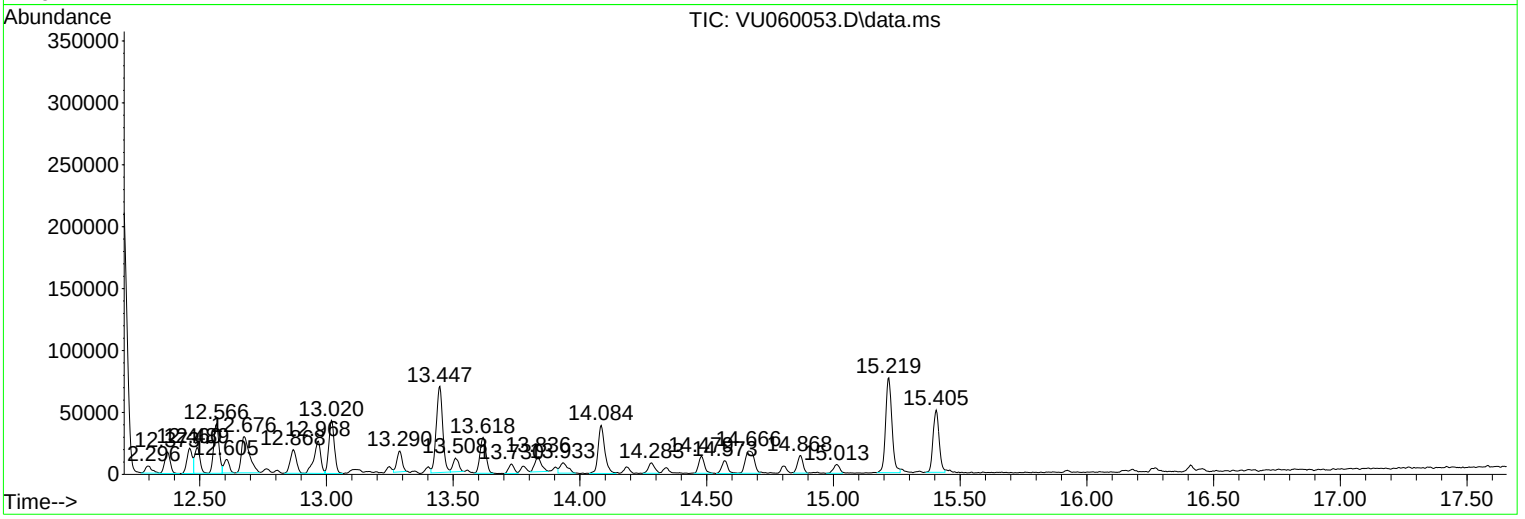
Instrument :
MSVOA_U
ClientSampleId :
YDZC5DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR071824WMA.M
Quant Title : TRACE VOA SFAM1.0

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







Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060053.D
Acq On : 19 Jul 2024 11:50
Operator : MD/SY
Sample : P3217-03DL 100X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZC5DL

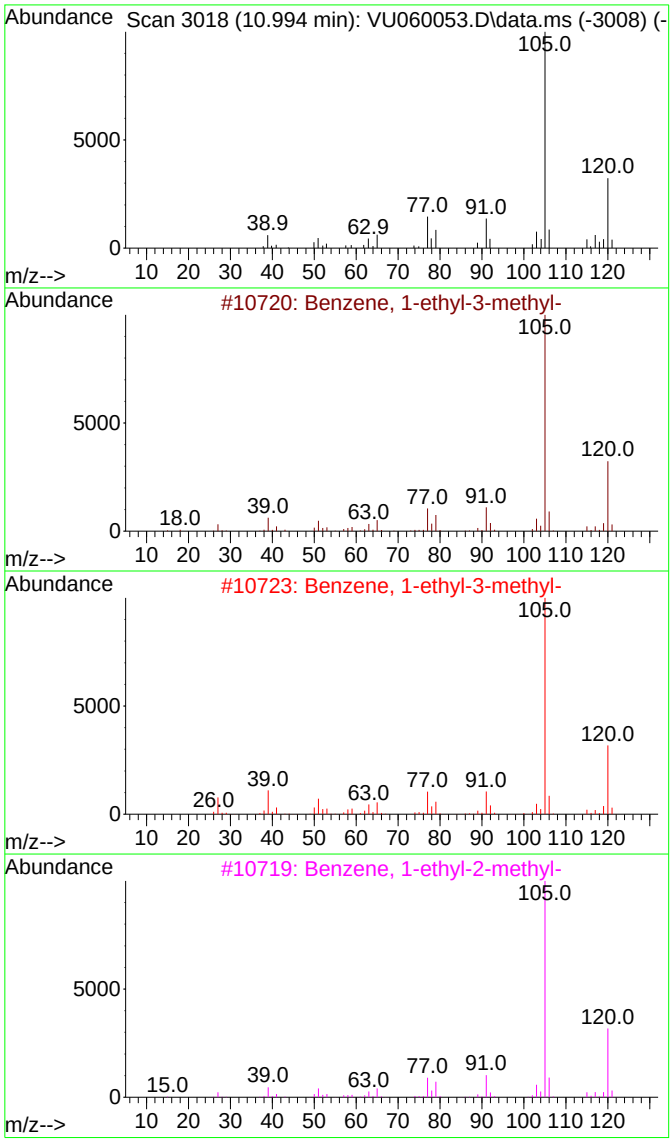
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR071824WMA.M
Quant Title : TRACE VOA SFAM1.0

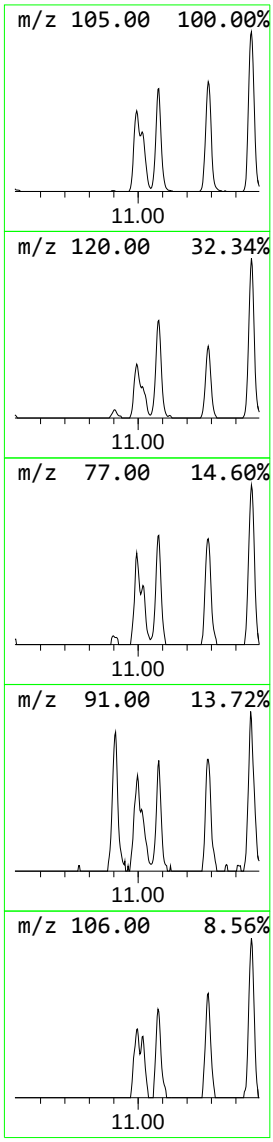
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Benzene, 1-ethyl-3-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.994	0.60 ug/L	56757	1,4-Dichlorobenzene-d4	11.807

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





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060053.D
Acq On : 19 Jul 2024 11:50
Operator : MD/SY
Sample : P3217-03DL 100X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZC5DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR071824WMA.M
Quant Title : TRACE VOA SFAM1.0

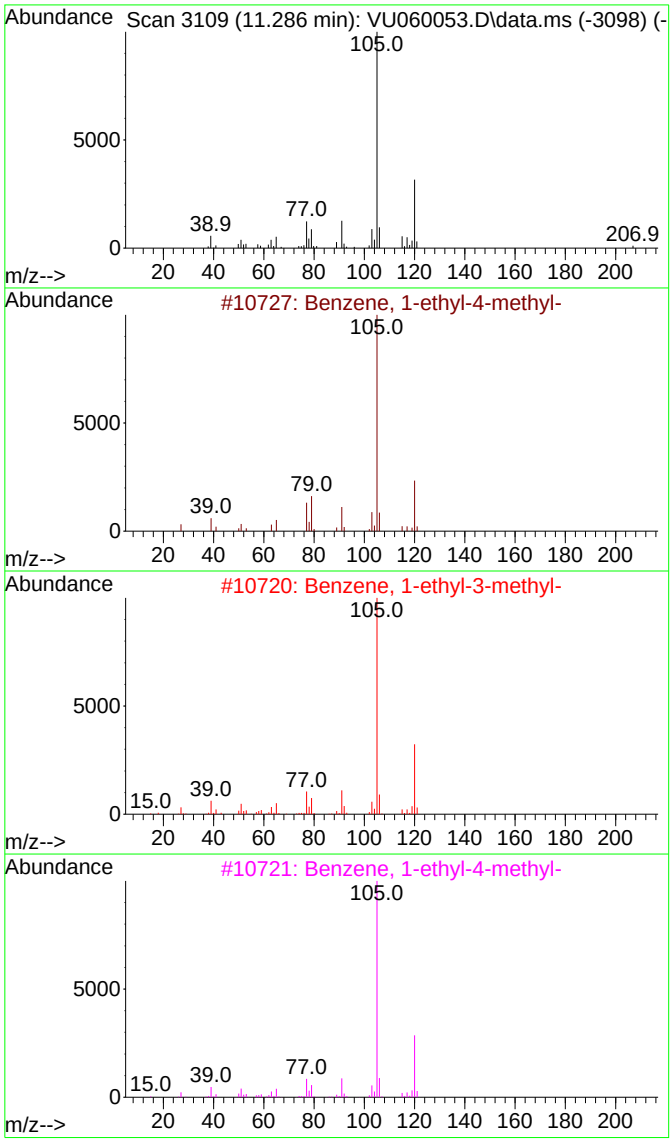
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

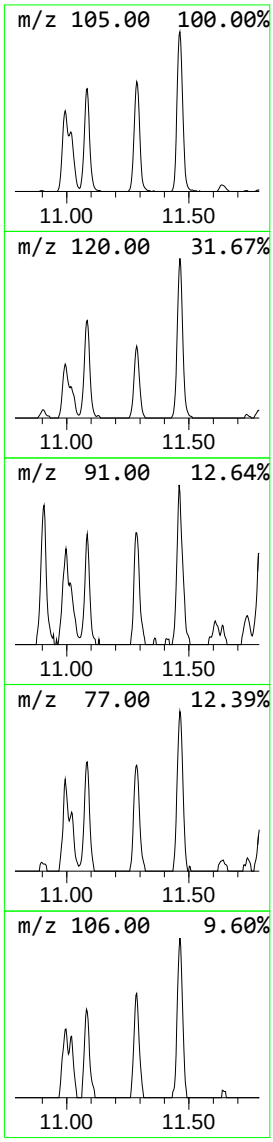
****

Peak Number 3 Benzene, 1-ethyl-4-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.287	0.78 ug/L	73336	1,4-Dichlorobenzene-d4	11.807

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





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060053.D
Acq On : 19 Jul 2024 11:50
Operator : MD/SY
Sample : P3217-03DL 100X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZC5DL

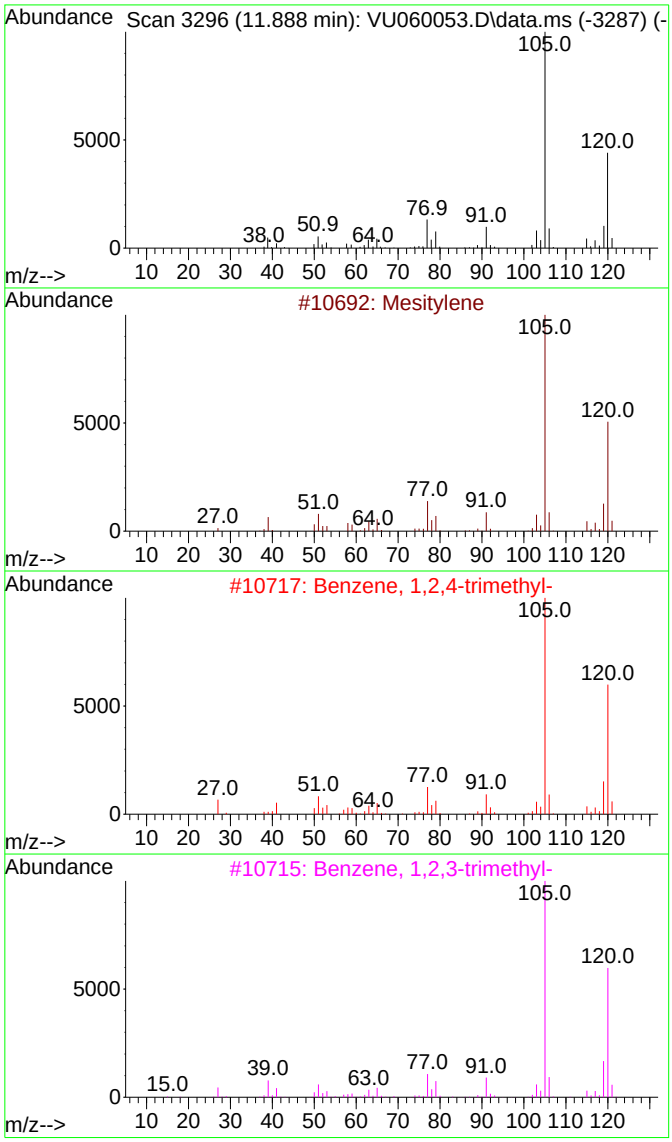
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR071824WMA.M
Quant Title : TRACE VOA SFAM1.0

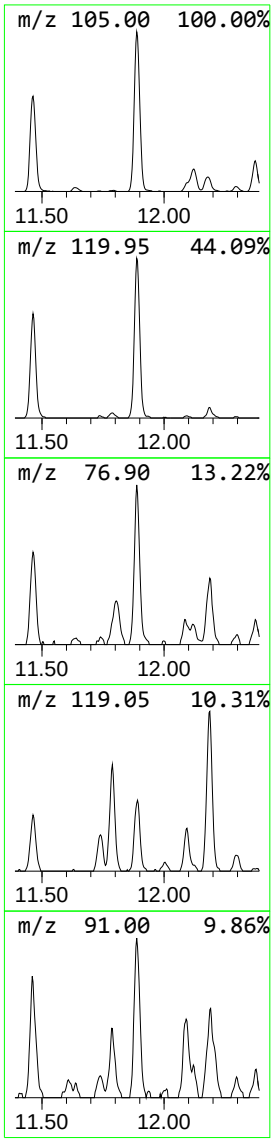
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.888	2.02 ug/L	189499	1,4-Dichlorobenzene-d4	11.807

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	97
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5	Mesitylene	120	C9H12	000108-67-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060053.D
Acq On : 19 Jul 2024 11:50
Operator : MD/SY
Sample : P3217-03DL 100X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZC5DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR071824WMA.M
Quant Title : TRACE VOA SFAM1.0

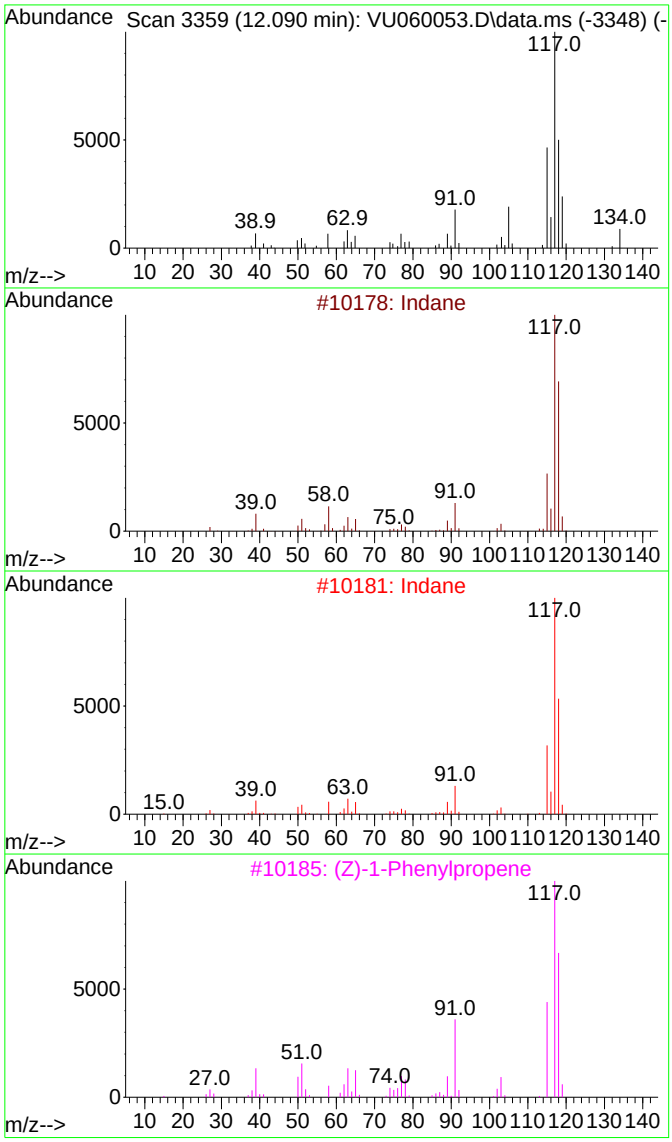
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

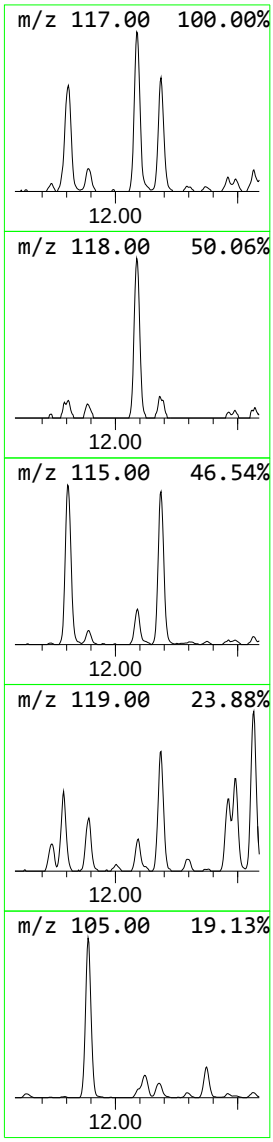
****

Peak Number 5 Indane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.090	0.82 ug/L	76679	1,4-Dichlorobenzene-d4	11.807

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	58
2	Indane	118	C9H10	000496-11-7	53
3	(Z)-1-Phenylpropene	118	C9H10	000766-90-5	53
4	Benzene, 2-propenyl-	118	C9H10	000300-57-2	49
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	47





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060053.D
Acq On : 19 Jul 2024 11:50
Operator : MD/SY
Sample : P3217-03DL 100X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZC5DL

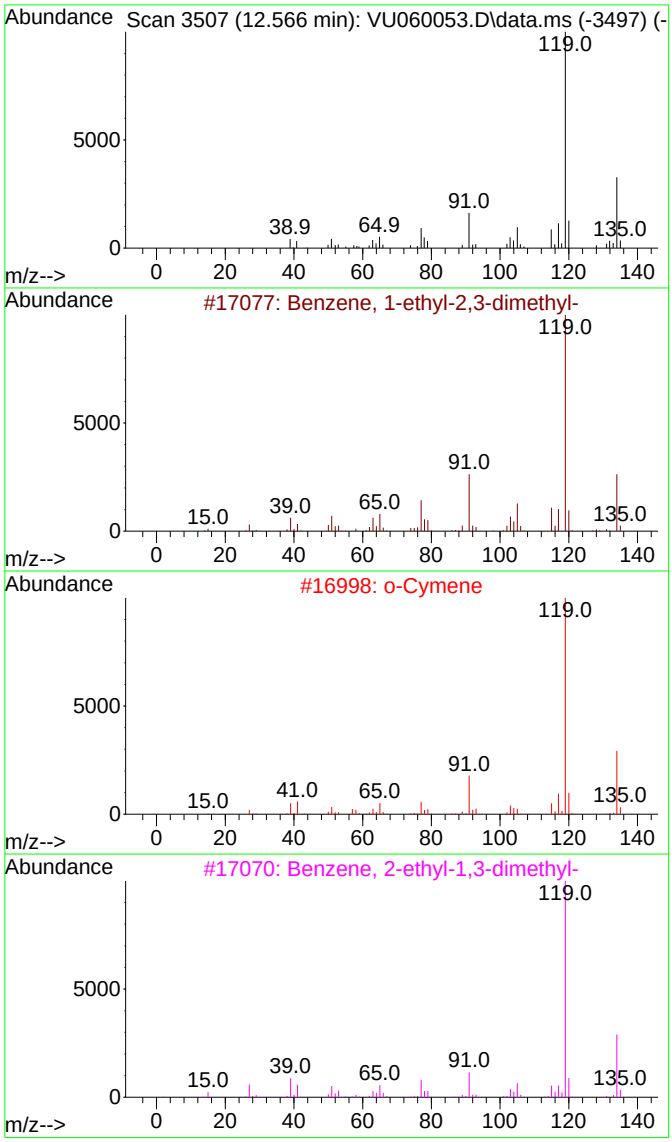
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR071824WMA.M
Quant Title : TRACE VOA SFAM1.0

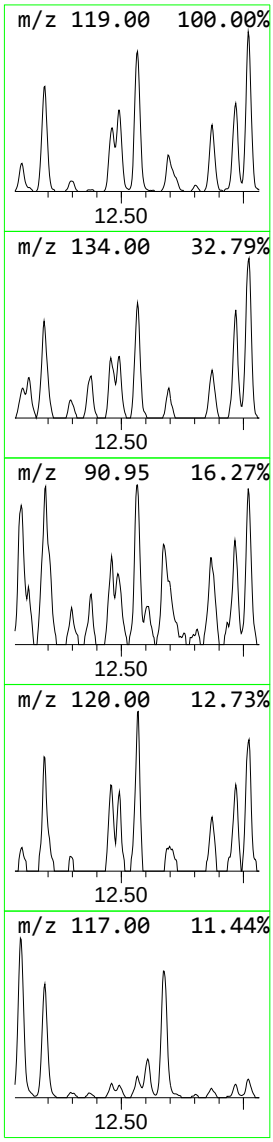
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.566	0.64 ug/L	60276	1,4-Dichlorobenzene-d4	11.807

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
2	o-Cymene	134	C10H14	000527-84-4	95
3	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060053.D
Acq On : 19 Jul 2024 11:50
Operator : MD/SY
Sample : P3217-03DL 100X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZC5DL

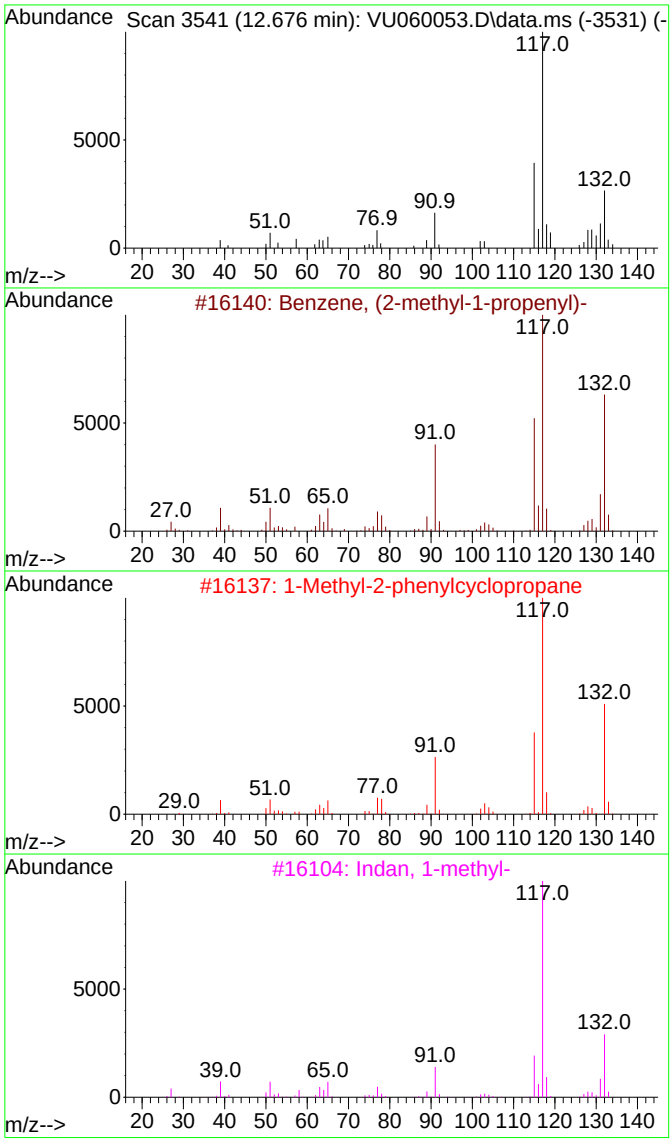
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR071824WMA.M
Quant Title : TRACE VOA SFAM1.0

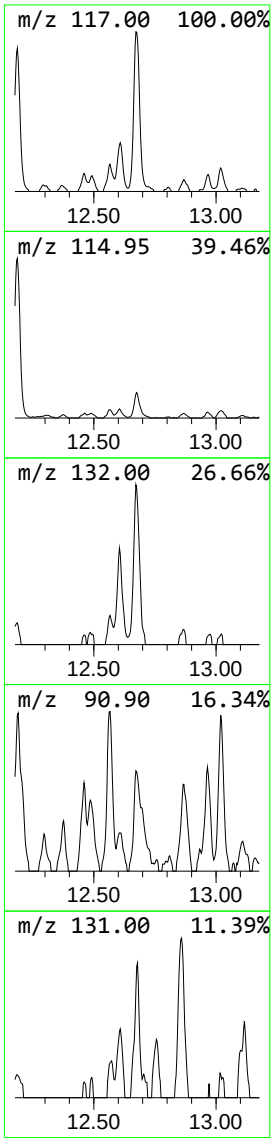
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, (2-methyl-1-propen... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.676	0.70 ug/L	65327	1,4-Dichlorobenzene-d4	11.807

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	86
2	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	83
3	Indan, 1-methyl-	132	C10H12	000767-58-8	76
4	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	72
5	Indan, 1-methyl-	132	C10H12	000767-58-8	70





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060053.D
Acq On : 19 Jul 2024 11:50
Operator : MD/SY
Sample : P3217-03DL 100X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZC5DL

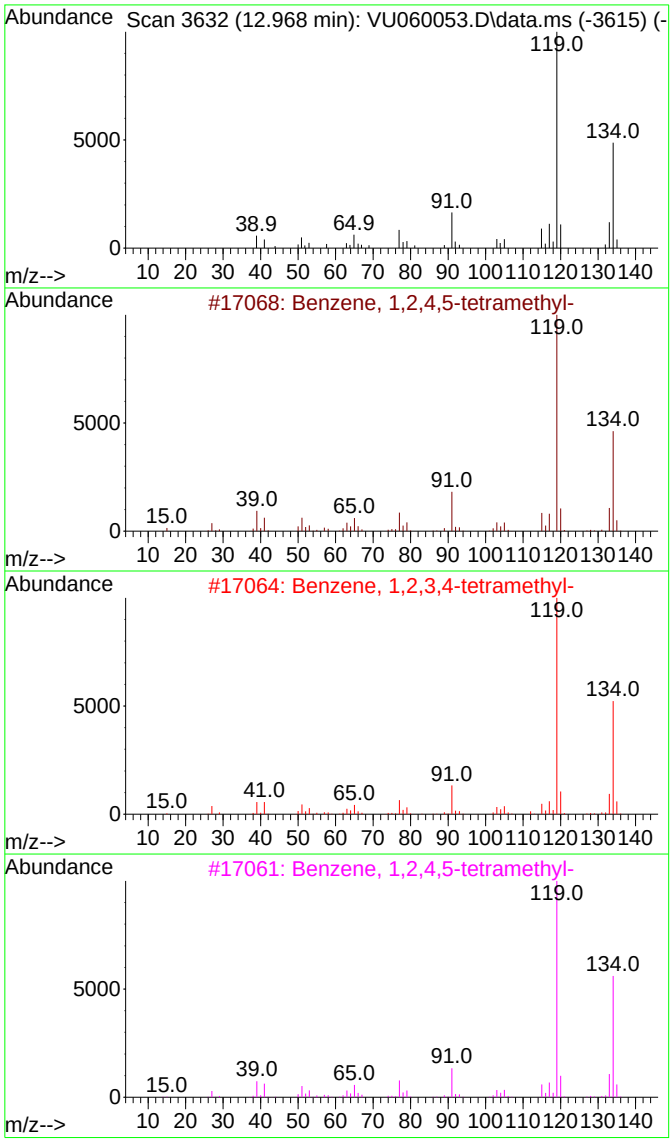
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR071824WMA.M
Quant Title : TRACE VOA SFAM1.0

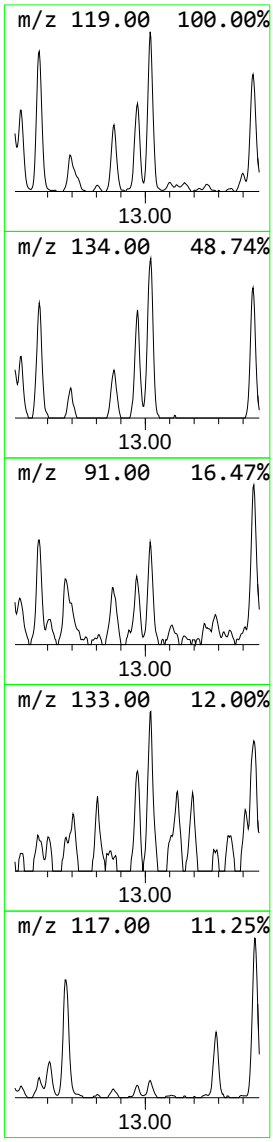
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.968	0.52 ug/L	48775	1,4-Dichlorobenzene-d4	11.807

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
4	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	91
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91





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Data File : VU060053.D
Acq On : 19 Jul 2024 11:50
Operator : MD/SY
Sample : P3217-03DL 100X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZC5DL

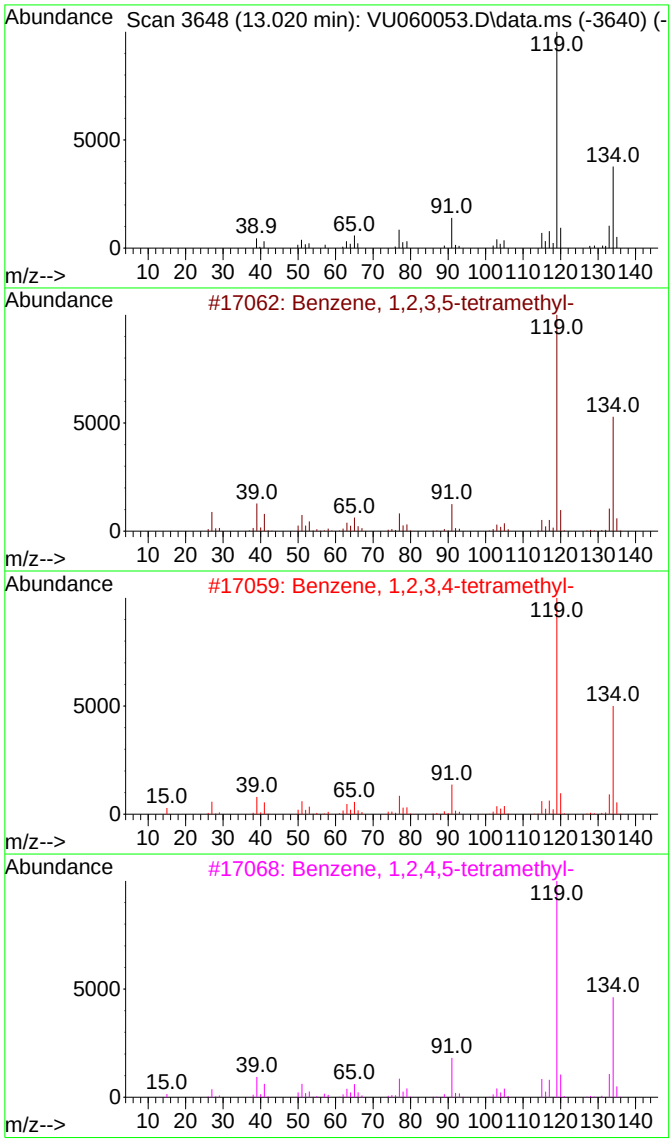
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR071824WMA.M
Quant Title : TRACE VOA SFAM1.0

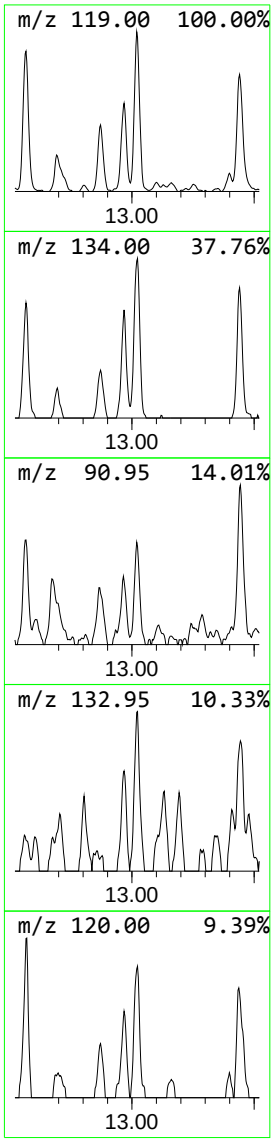
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.020	0.69 ug/L	65144	1,4-Dichlorobenzene-d4	11.807

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
4	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060053.D
Acq On : 19 Jul 2024 11:50
Operator : MD/SY
Sample : P3217-03DL 100X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZC5DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR071824WMA.M
Quant Title : TRACE VOA SFAM1.0

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

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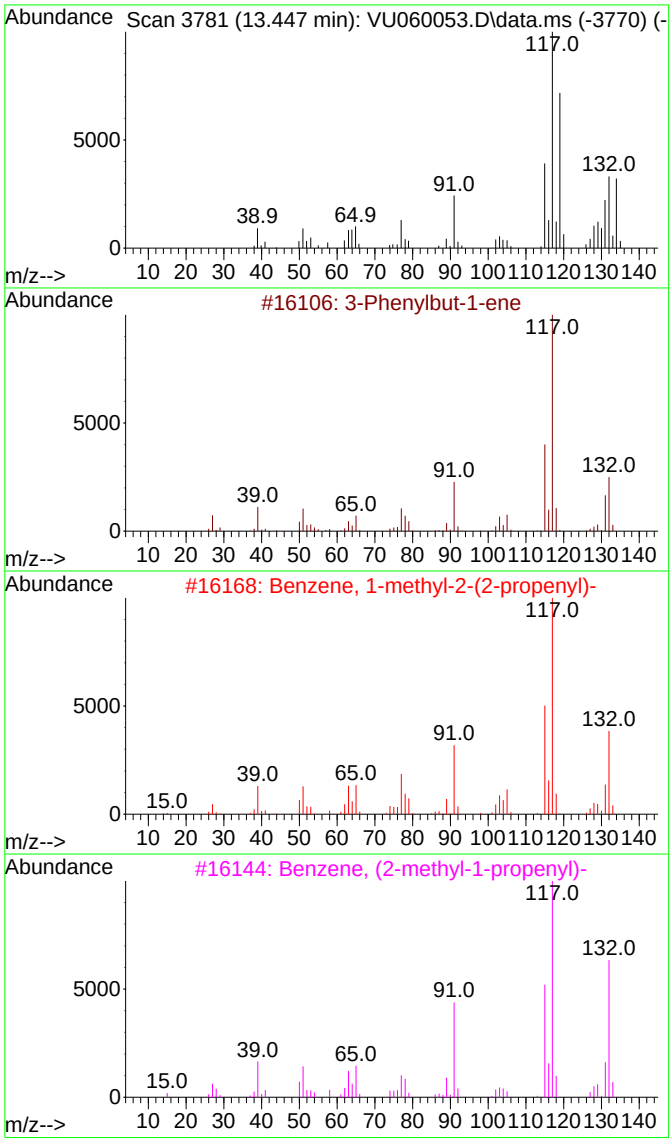
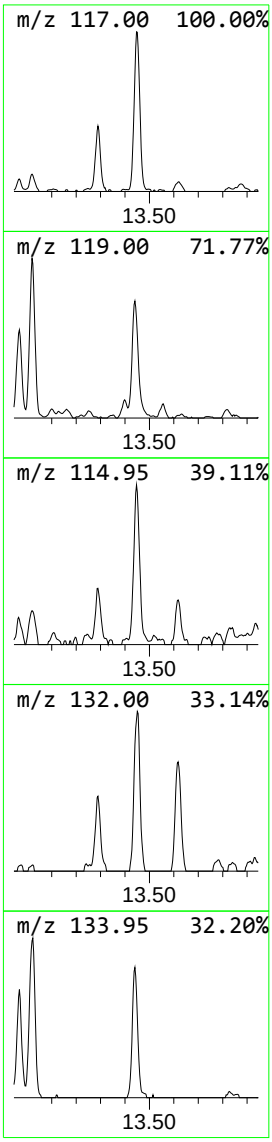
Peak Number 10 3-Phenylbut-1-ene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.447	1.31 ug/L	123278	1,4-Dichlorobenzene-d4	11.807

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Phenylbut-1-ene	132	C10H12	000934-10-1	64
2	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	60
3	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	60
4	Benzene, 2-butenyl-	132	C10H12	001560-06-1	60
5	Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	60





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060053.D
Acq On : 19 Jul 2024 11:50
Operator : MD/SY
Sample : P3217-03DL 100X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZC5DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR071824WMA.M
Quant Title : TRACE VOA SFAM1.0

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

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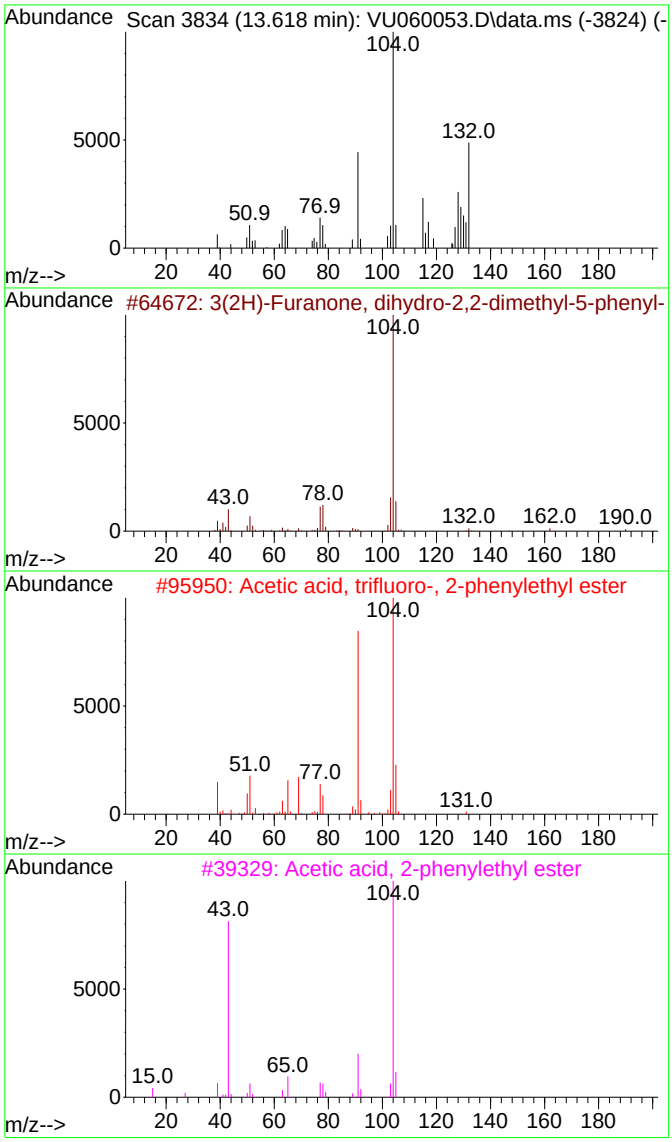
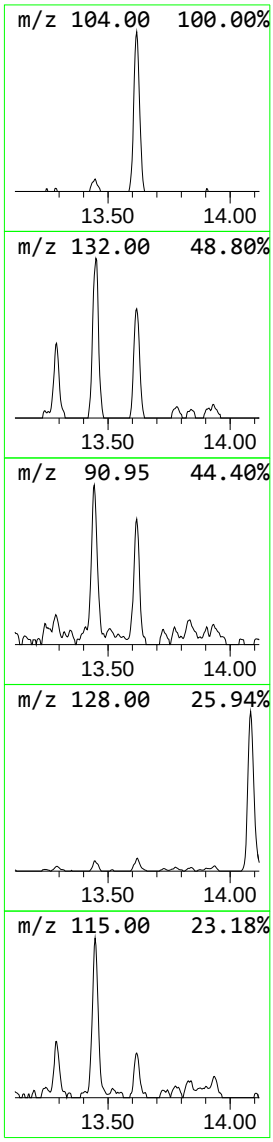
Peak Number 11 unknown-01 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.618	0.51 ug/L	48189	1,4-Dichlorobenzene-d4	11.807

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3(2H)-Furanone, dihydro-2,2-dime...	190	C12H14O2	063678-00-2	27
2	Acetic acid, trifluoro-, 2-pheny...	218	C10H9F3O2	055419-66-4	27
3	Acetic acid, 2-phenylethyl ester	164	C10H12O2	000103-45-7	27
4	Acetic acid, trichloro-, 2-pheny...	266	C10H9Cl3O2	071965-07-6	27
5	2-Phenylethylamine, N,N-di(trifl...	313	C12H9F6NO2	1000463-67-7	27





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060053.D
Acq On : 19 Jul 2024 11:50
Operator : MD/SY
Sample : P3217-03DL 100X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZC5DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR071824WMA.M
Quant Title : TRACE VOA SFAM1.0

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

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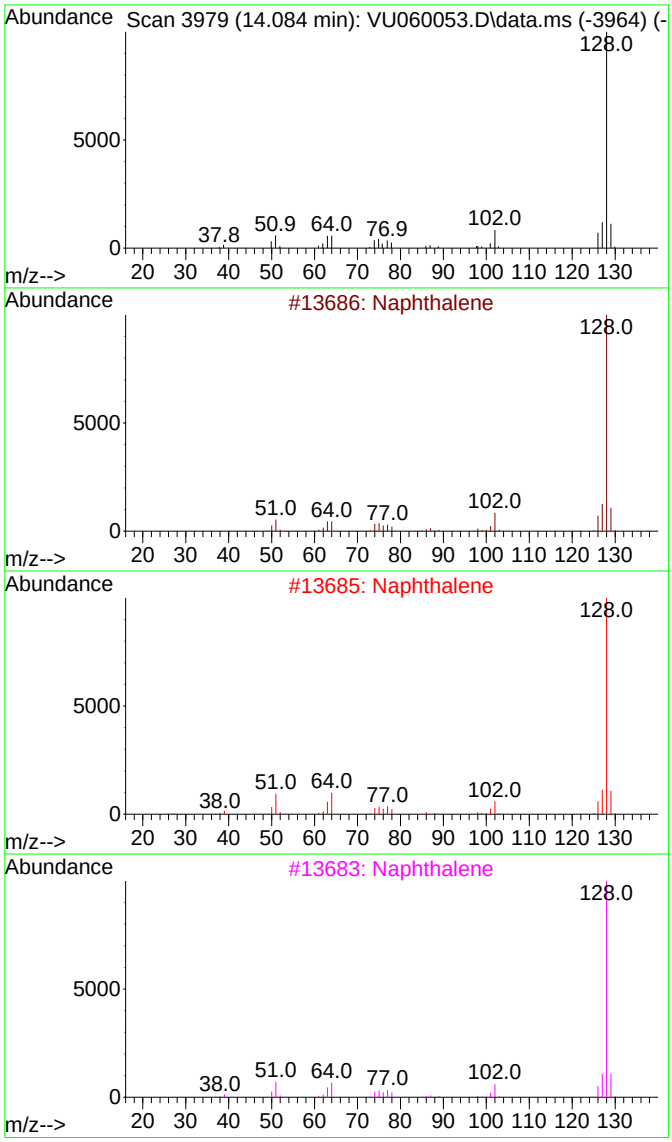
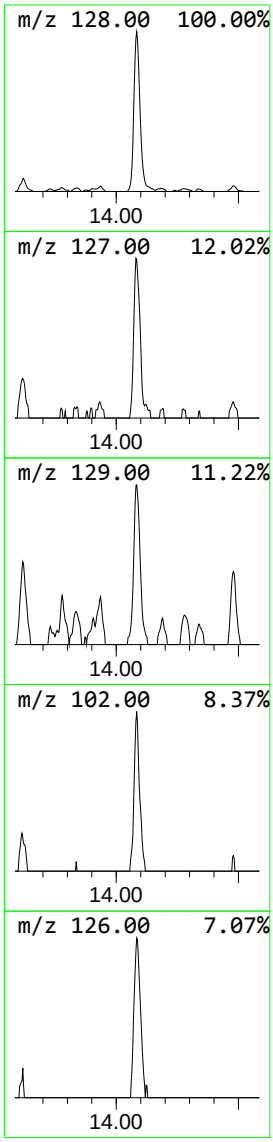
Peak Number 12 Azulene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.084	0.75 ug/L	70592	1,4-Dichlorobenzene-d4	11.807

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene	128	C10H8	000091-20-3	95
2	Naphthalene	128	C10H8	000091-20-3	91
3	Naphthalene	128	C10H8	000091-20-3	91
4	Azulene	128	C10H8	000275-51-4	91
5	Azulene	128	C10H8	000275-51-4	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU071824\
Data File : VU060053.D
Acq On : 19 Jul 2024 11:50
Operator : MD/SY
Sample : P3217-03DL 100X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 48 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
YDZC5DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR071824WMA.M
Quant Title : TRACE VOA SFAM1.0

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-ethy...	10.994	0.6	ug/L	56757	3	11.807	469565	5.0
Benzene, 1-ethy...	11.287	0.8	ug/L	73336	3	11.807	469565	5.0
Benzene, 1,2,3-...	11.888	2.0	ug/L	189499	3	11.807	469565	5.0
Indane	12.090	0.8	ug/L	76679	3	11.807	469565	5.0
Benzene, 1-ethy...	12.566	0.6	ug/L	60276	3	11.807	469565	5.0
Benzene, (2-met...	12.676	0.7	ug/L	65327	3	11.807	469565	5.0
Benzene, 1,2,4,...	12.968	0.5	ug/L	48775	3	11.807	469565	5.0
Benzene, 1,2,3,...	13.020	0.7	ug/L	65144	3	11.807	469565	5.0
3-Phenylbut-1-ene	13.447	1.3	ug/L	123278	3	11.807	469565	5.0
unknown-01	13.618	0.5	ug/L	48189	3	11.807	469565	5.0
Azulene	14.084	0.8	ug/L	70592	3	11.807	469565	5.0