LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN021424\
Data File : VN081032.D
Acq On : 14 Feb 2024 18:07
Operator : JC\MD
Sample : P1459-03 20X
Misc : 1.01g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WAL-OILY-DEBRIS

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_N\methods\82N021424W.M
Title : SW846 8260

Signal : TIC: VN081032.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.659	114	120	129	rBV	54798	104119	3.65%	0.246%
2	2.818	139	147	148	rBV3	39472	81194	2.85%	0.192%
3	8.165	1046	1056	1061	rBV	385291	931101	32.67%	2.200%
4	8.224	1061	1066	1077	rVB	779281	1716383	60.23%	4.055%
5	8.577	1117	1126	1137	rBV	410925	948252	33.28%	2.240%
6	9.100	1204	1215	1228	rBV	1129075	2373491	83.29%	5.607%
7	9.335	1247	1255	1264	rBV	117928	266708	9.36%	0.630%
8	9.571	1288	1295	1306	rBV3	49781	144636	5.08%	0.342%
9	10.494	1447	1452	1457	rBV5	27795	49758	1.75%	0.118%
10	10.571	1457	1465	1471	rVV	1448680	2708799	95.05%	6.400%
11	10.629	1471	1475	1485	rVB	97660	182981	6.42%	0.432%
12	10.747	1492	1495	1502	rVB7	22583	37363	1.31%	0.088%
13	10.912	1518	1523	1532	rVB5	31602	67865	2.38%	0.160%
14	11.000	1534	1538	1545	rBV5	34830	70169	2.46%	0.166%
15	11.276	1581	1585	1589	rBV6	20369	33264	1.17%	0.079%
16	11.418	1602	1609	1615	rVV3	72673	166316	5.84%	0.393%
17	11.470	1615	1618	1624	rVB5	29803	52621	1.85%	0.124%
18	11.647	1644	1648	1650	rBV4	23051	37796	1.33%	0.089%
19	11.682	1650	1654	1657	rVV6	25973	44338	1.56%	0.105%
20	11.747	1662	1665	1671	rVB5	22796	42438	1.49%	0.100%
21	11.865	1678	1685	1696	rVB	1570364	2849730	100.00%	6.733%
22	11.965	1697	1702	1707	rBV	115821	203310	7.13%	0.480%
23	12.070	1713	1720	1724	rBV	347104	674741	23.68%	1.594%
24	12.153	1732	1734	1740	rVB3	52770	76550	2.69%	0.181%
25	12.400	1765	1776	1781	rBV3	218321	596422	20.93%	1.409%
26	12.535	1796	1799	1803	rBV2	80273	123764	4.34%	0.292%
27	12.623	1810	1814	1818	rBV2	101766	154316	5.42%	0.365%
28	12.794	1841	1843	1847	rBV5	32417	55674	1.95%	0.132%
29	12.853	1847	1853	1860	rVB	1116782	1936621	67.96%	4.575%
30	13.035	1878	1884	1889	rBV3	93835	162284	5.69%	0.383%
31	13.100	1889	1895	1899	rBV	266909	518620	18.20%	1.225%
32	13.176	1903	1908	1914	rVV4	329988	668252	23.45%	1.579%
33	13.229	1914	1917	1922	rVB4	65232	102561	3.60%	0.242%
34	13.488	1955	1961	1967	rBV	513597	941731	33.05%	2.225%
35	13.617	1979	1983	1987	rBV4	99748	178075	6.25%	0.421%
36	13.682	1987	1994	1998	rVV5	223240	526289	18.47%	1.243%

LSC Area Percent Report

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Operator : JC\MD
Sample : P1459-03 20X
Misc : 1.01g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WAL-OILY-DEBRIS

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_N\methods\82N021424W.M
Title : SW846 8260

37	13.794	2007	2013	2023	rVB2	1402498	2749087	96.47%	6.495%
38	13.988	2041	2046	2049	rBV2	216067	364086	12.78%	0.860%
39	14.123	2063	2069	2078	rBV4	602810	1320710	46.35%	3.120%
40	14.247	2083	2090	2093	rBV6	122517	311152	10.92%	0.735%
41	14.335	2101	2105	2110	rVB2	242430	385535	13.53%	0.911%
42	14.441	2119	2123	2129	rVB2	345129	537479	18.86%	1.270%
43	14.506	2130	2134	2140	rBV5	291656	569530	19.99%	1.346%
44	14.576	2142	2146	2150	rVV5	335902	536631	18.83%	1.268%
45	14.629	2150	2155	2158	rVV	713121	1171967	41.13%	2.769%
46	14.664	2158	2161	2169	rVB2	502893	1005166	35.27%	2.375%
47	14.794	2176	2183	2188	rBV4	569091	1197435	42.02%	2.829%
48	14.847	2188	2192	2202	rVB9	397845	843793	29.61%	1.993%
49	15.058	2220	2228	2233	rVV6	514166	1473068	51.69%	3.480%
50	15.111	2233	2237	2245	rVB5	554667	968894	34.00%	2.289%
51	15.217	2251	2255	2260	rBV2	586311	1162981	40.81%	2.748%
52	15.270	2260	2264	2268	rVV3	530072	864294	30.33%	2.042%
53	15.323	2269	2273	2282	rVB7	330948	756817	26.56%	1.788%
54	15.700	2330	2337	2344	rBV8	774781	2189828	76.84%	5.174%
55	15.929	2372	2376	2379	rBV4	251441	489574	17.18%	1.157%
56	16.170	2412	2417	2423	rBV5	866433	1838651	64.52%	4.344%
57	16.711	2504	2509	2516	rBV2	846268	1762339	61.84%	4.164%

Sum of corrected areas: 42327549

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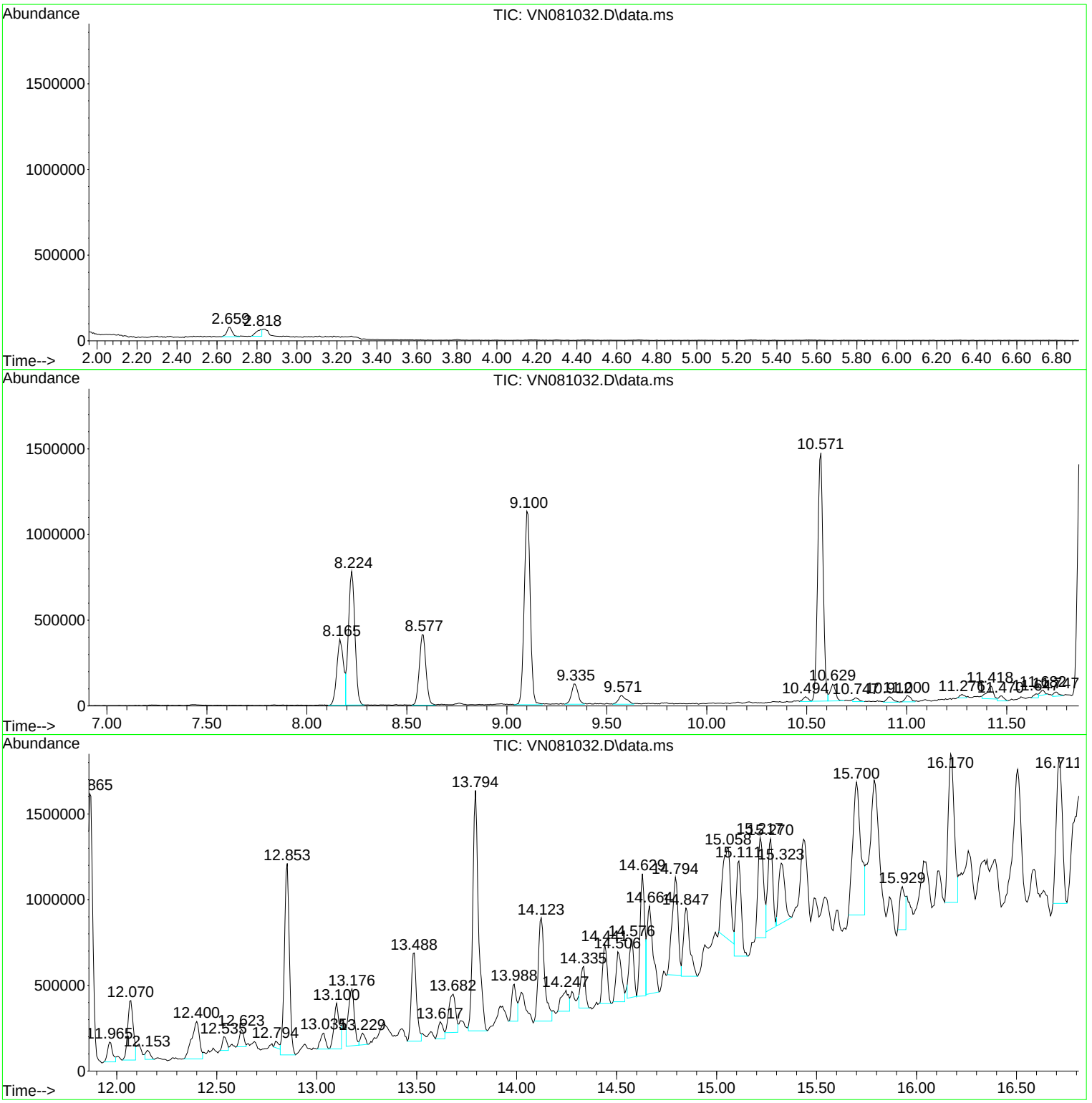
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WAL-OILY-DEBRIS

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021424W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN021424\
Data File : VN081032.D
Acq On : 14 Feb 2024 18:07
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Sample : P1459-03 20X
Misc : 1.01g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WAL-OILY-DEBRIS

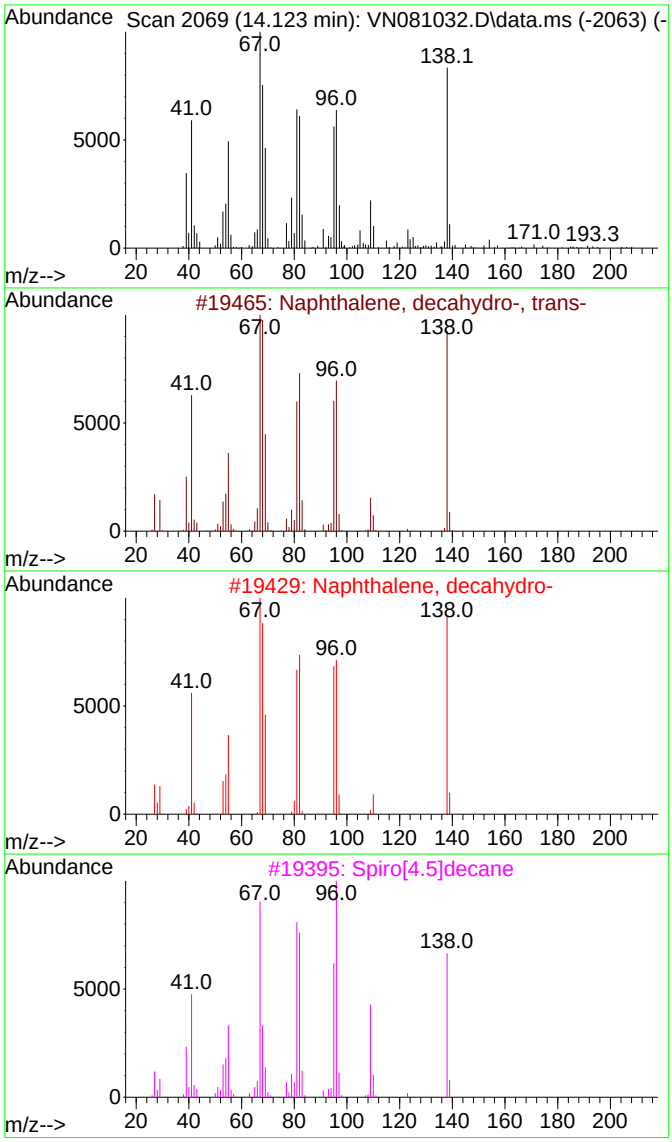
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021424W.M
Quant Title : SW846 8260

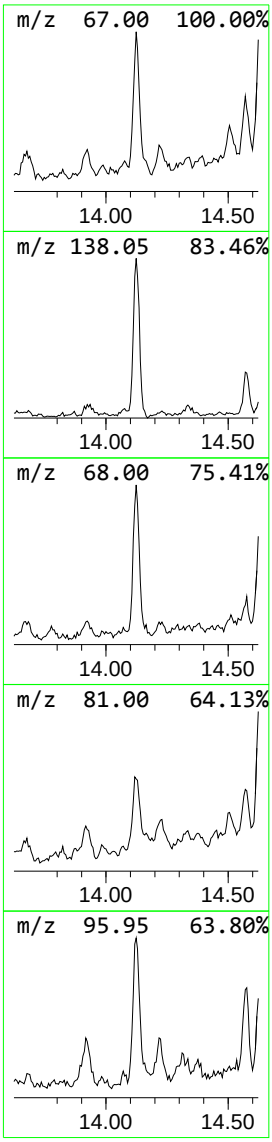
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Naphthalene, decahydro-, tr... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.123	24.02 ug/l	1320710	1,4-Dichlorobenzene-d4	13.794

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	97
2	Naphthalene, decahydro-	138	C10H18	000091-17-8	97
3	Spiro[4.5]decane	138	C10H18	000176-63-6	76
4	Naphthalene, decahydro-, cis-	138	C10H18	000493-01-6	76
5	2-Cyclohexen-1-one, 4-(1-methyle...	138	C9H14O	000500-02-7	74





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN021424\
Data File : VN081032.D
Acq On : 14 Feb 2024 18:07
Operator : JC\MD
Sample : P1459-03 20X
Misc : 1.01g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WAL-OILY-DEBRIS

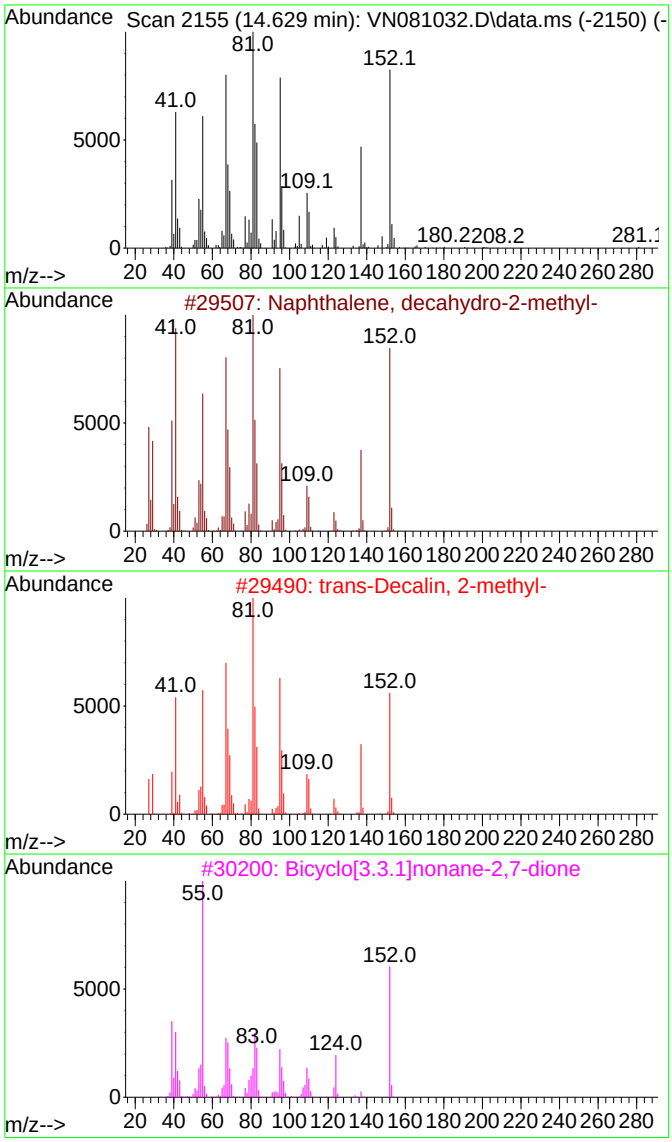
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021424W.M
Quant Title : SW846 8260

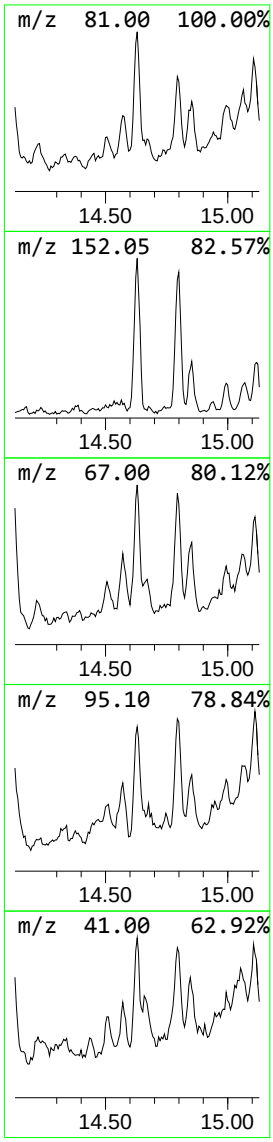
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Naphthalene, decahydro-2-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.629	21.32 ug/l	1171970	1,4-Dichlorobenzene-d4	13.794

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	97
2	trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	96
3	Bicyclo[3.3.1]nonane-2,7-dione	152	C9H12O2	1000210-30-3	70
4	Cyclodecene, 1-methyl-	152	C11H20	066633-38-3	68
5	Decalin, syn-1-methyl-, cis-	152	C11H20	1000158-89-1	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN021424\
Data File : VN081032.D
Acq On : 14 Feb 2024 18:07
Operator : JC\MD
Sample : P1459-03 20X
Misc : 1.01g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WAL-OILY-DEBRIS

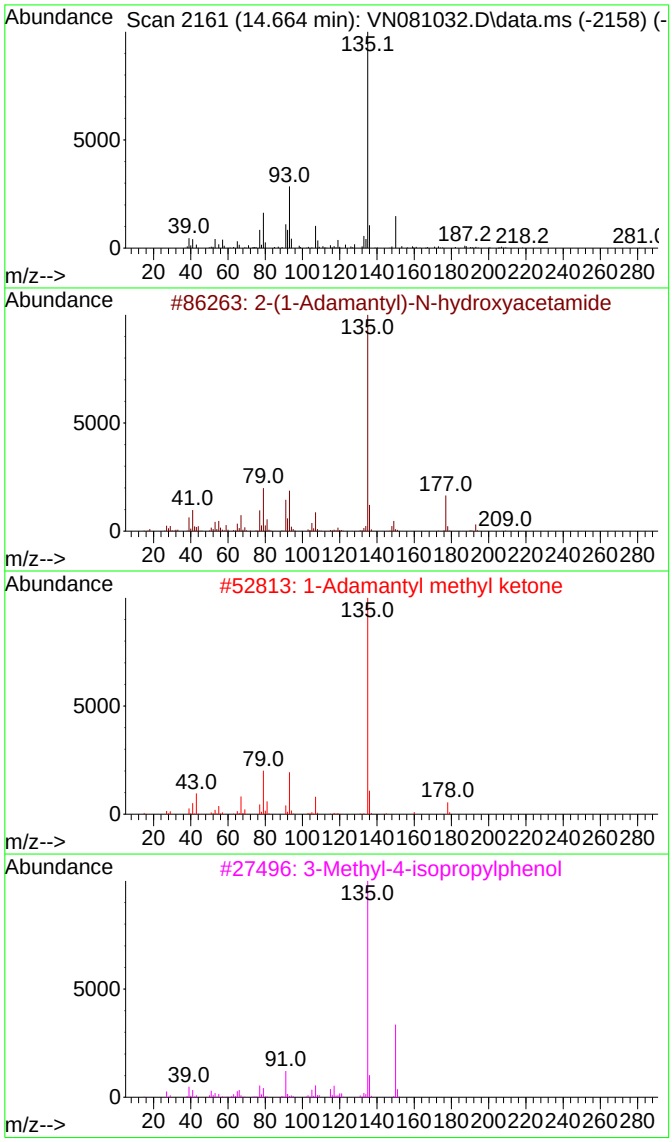
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021424W.M
Quant Title : SW846 8260

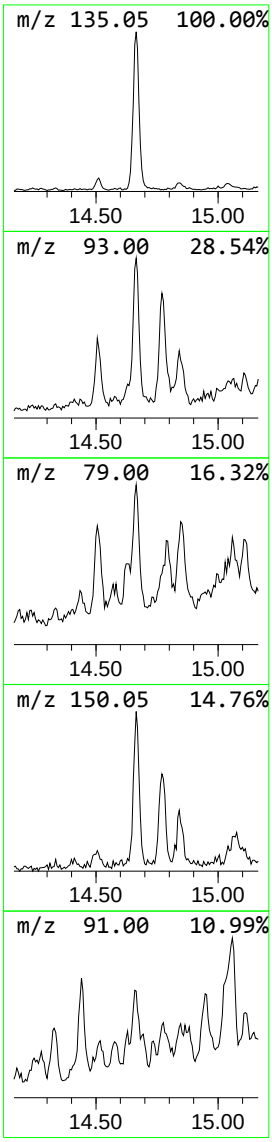
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 2-(1-Adamantyl)-N-hydroxyac... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.664	18.28 ug/l	1005170	1,4-Dichlorobenzene-d4	13.794

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-(1-Adamantyl)-N-hydroxyacetamide	209	C12H19NO2	136561-40-5	64
2	1-Adamantyl methyl ketone	178	C12H18O	001660-04-4	64
3	3-Methyl-4-isopropylphenol	150	C10H14O	003228-02-2	64
4	4-Cyanobenzoic acid, 2-(1-adaman...	309	C20H23NO2	1000292-45-1	64
5	1-Adamantanemethanol	166	C11H18O	000770-71-8	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN021424\
Data File : VN081032.D
Acq On : 14 Feb 2024 18:07
Operator : JC\MD
Sample : P1459-03 20X
Misc : 1.01g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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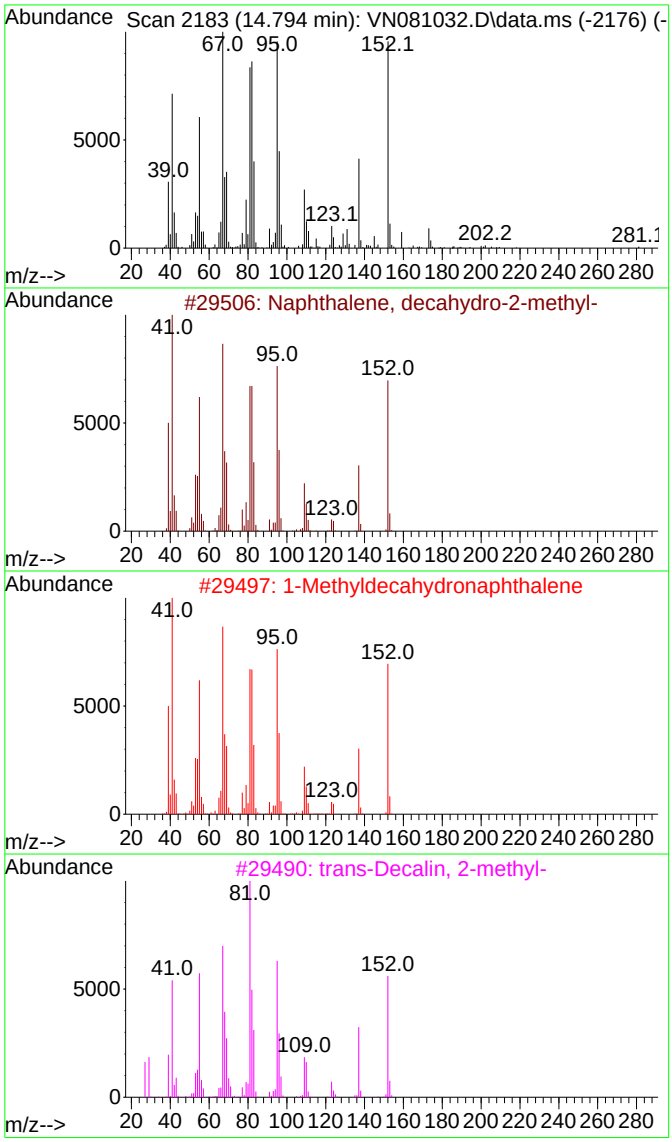
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021424W.M
Quant Title : SW846 8260

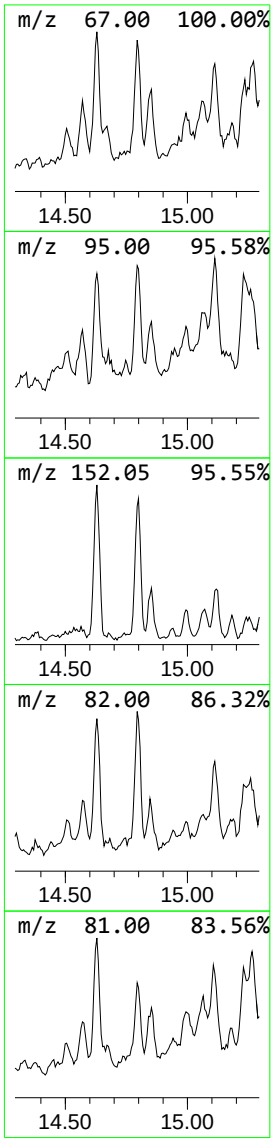
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 1-Methyldecahydronaphthalene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.794	21.78 ug/l	1197440	1,4-Dichlorobenzene-d4	13.794

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	90
2	1-Methyldecahydronaphthalene	152	C11H20	002958-75-0	90
3	trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	89
4	cis-Decalin, 2-syn-methyl-	152	C11H20	1000155-85-6	62
5	Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	62





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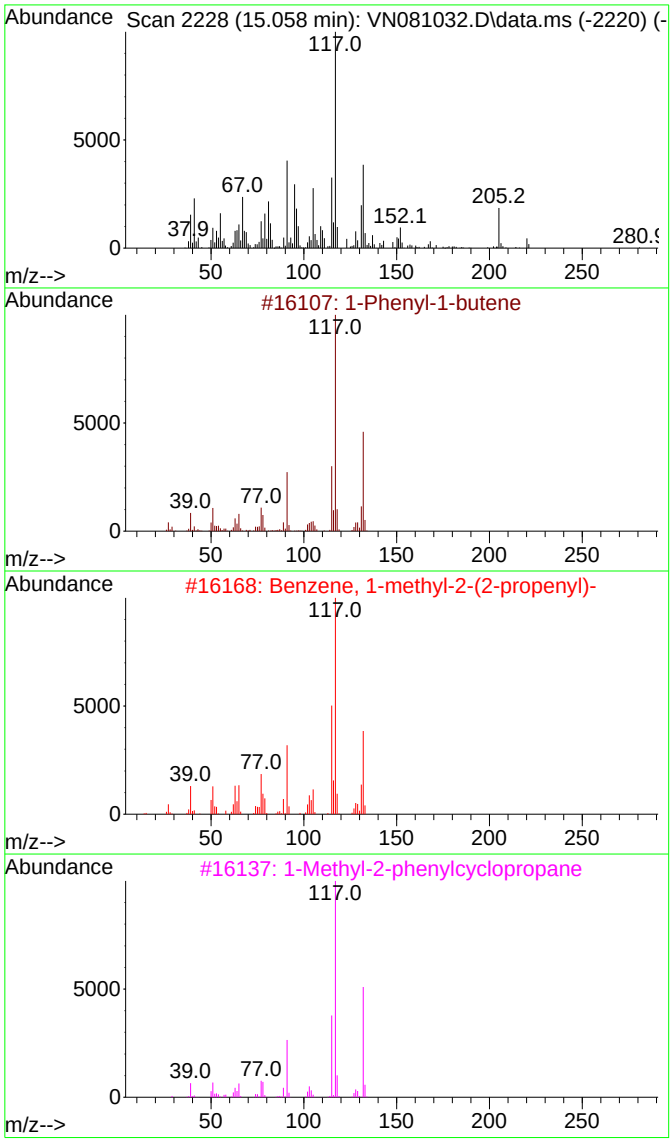
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021424W.M
Quant Title : SW846 8260

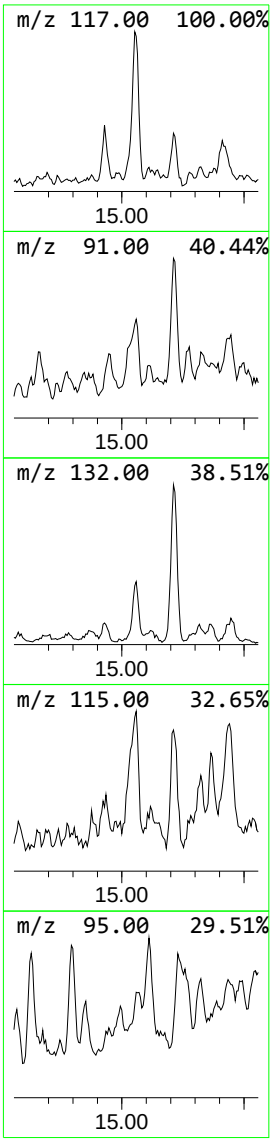
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 1-Phenyl-1-butene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.059	26.79 ug/l	1473070	1,4-Dichlorobenzene-d4	13.794

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenyl-1-butene	132	C10H12	000824-90-8	94
2	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	89
3	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	83
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
5	Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	76





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Misc : 1.01g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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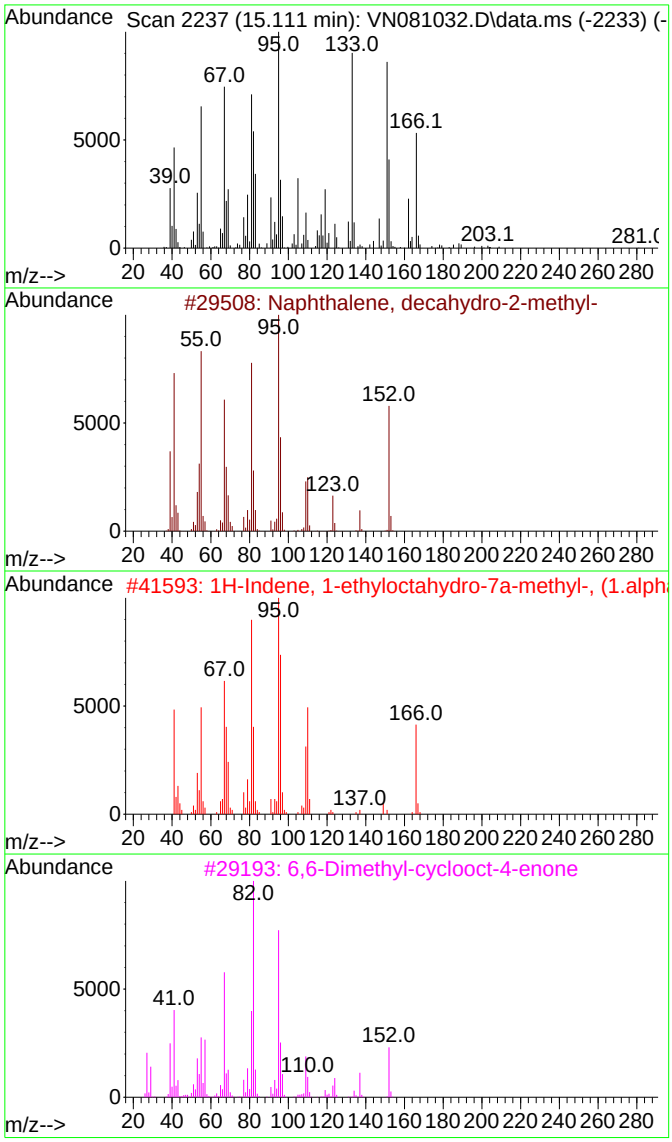
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021424W.M
Quant Title : SW846 8260

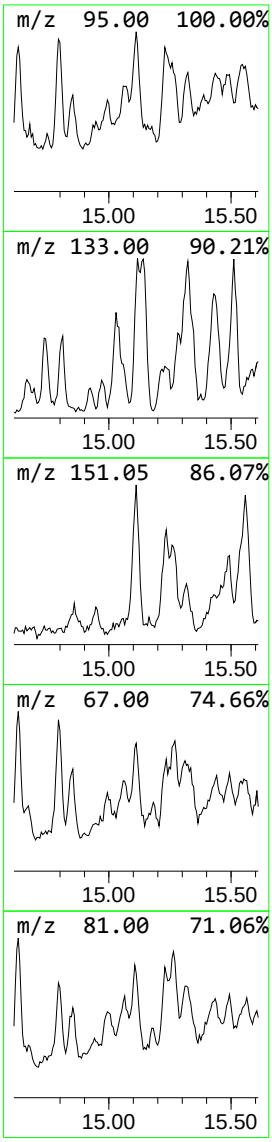
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 unknown15.111 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.111	17.62 ug/l	968894	1,4-Dichlorobenzene-d4	13.794

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	38
2	1H-Indene, 1-ethyloctahydro-7a-m...	166	C12H22	056324-71-1	38
3	6,6-Dimethyl-cyclooct-4-enone	152	C10H16O	1000194-10-0	30
4	Cycloheptanone, 2-(2-methylpropy...	166	C11H18O	120694-89-5	25
5	Decalin, anti-1-methyl-, cis-	152	C11H20	1000158-89-0	25





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Data File : VN081032.D
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Instrument :
MSVOA_N
ClientSampleId :
WAL-OILY-DEBRIS

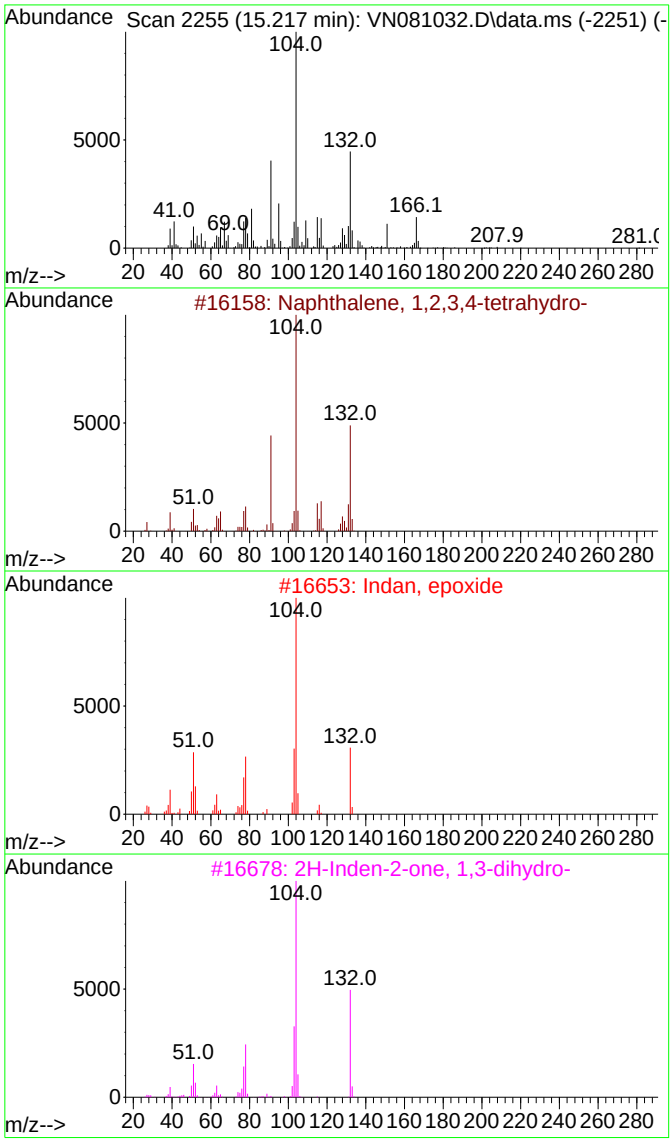
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021424W.M
Quant Title : SW846 8260

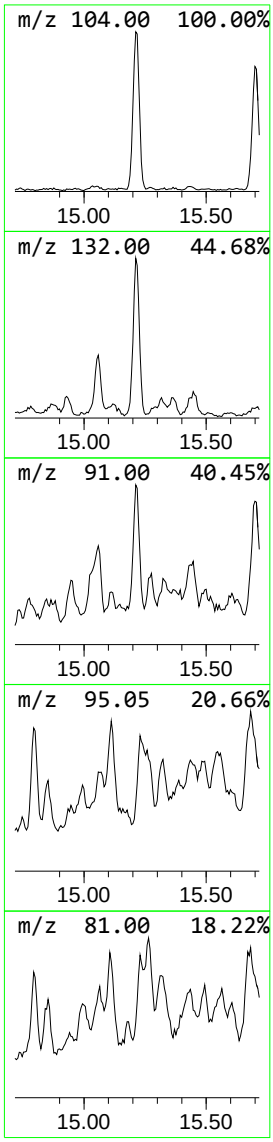
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.217	21.15 ug/l	1162980	1,4-Dichlorobenzene-d4	13.794

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	97
2	Indan, epoxide	132	C9H8O	000768-22-9	49
3	2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	46
4	Benzene, [2-(1,1-dimethylethyl)c...	174	C13H18	1000362-12-2	38
5	2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN021424\
Data File : VN081032.D
Acq On : 14 Feb 2024 18:07
Operator : JC\MD
Sample : P1459-03 20X
Misc : 1.01g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WAL-OILY-DEBRIS

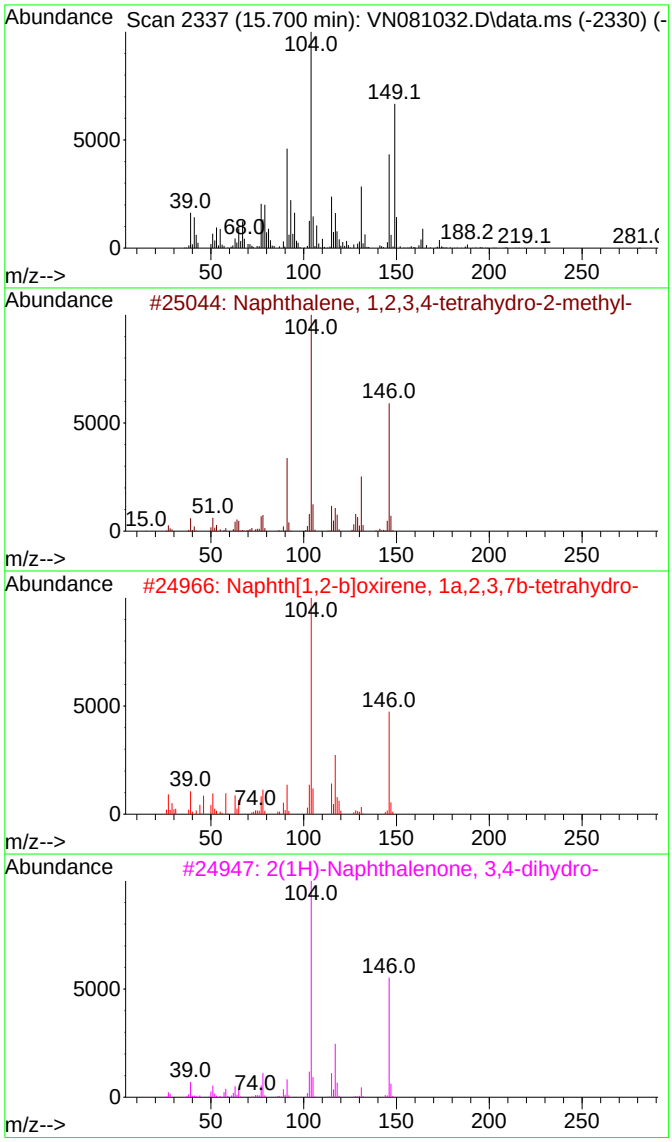
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021424W.M
Quant Title : SW846 8260

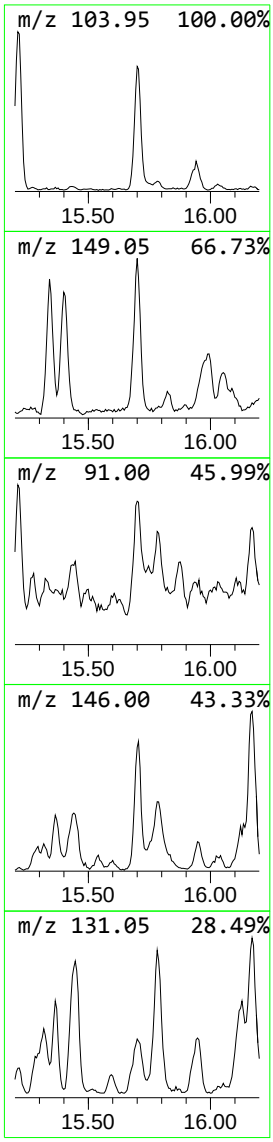
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.700	39.83 ug/l	2189830	1,4-Dichlorobenzene-d4	13.794

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	64
2	Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	46
3	2(1H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000530-93-8	30
4	Bromoacetic acid, 2-phenylethyl ...	242	C10H11BrO2	003785-33-9	22
5	Acetic acid, 2-phenylethyl ester	164	C10H12O2	000103-45-7	22





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN021424\
Data File : VN081032.D
Acq On : 14 Feb 2024 18:07
Operator : JC\MD
Sample : P1459-03 20X
Misc : 1.01g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WAL-OILY-DEBRIS

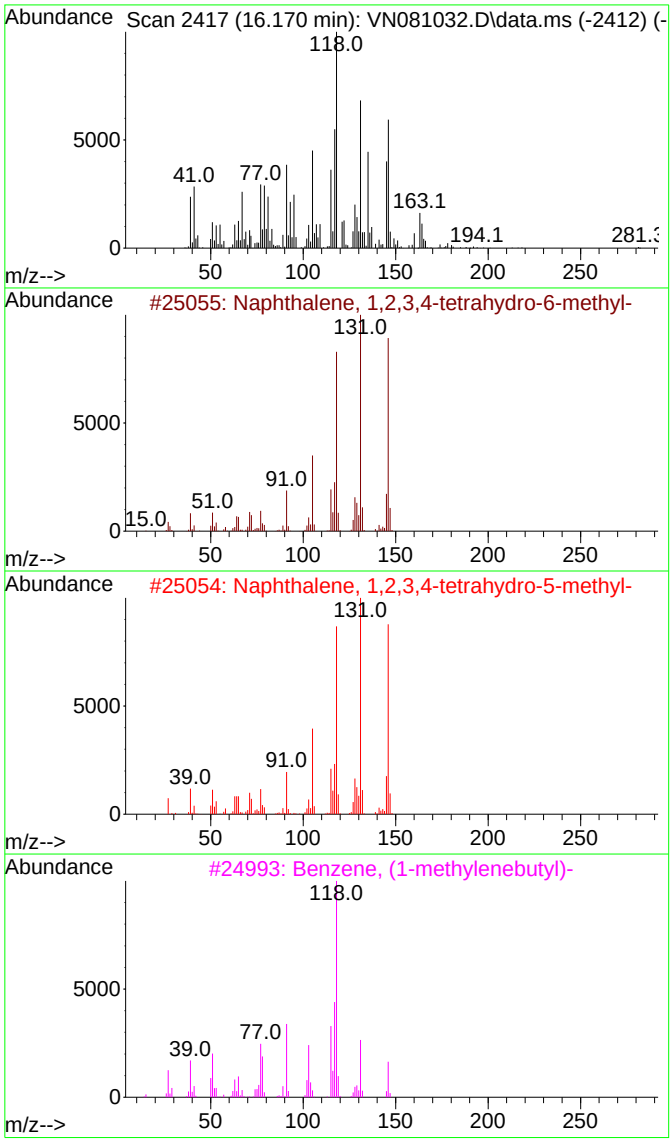
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021424W.M
Quant Title : SW846 8260

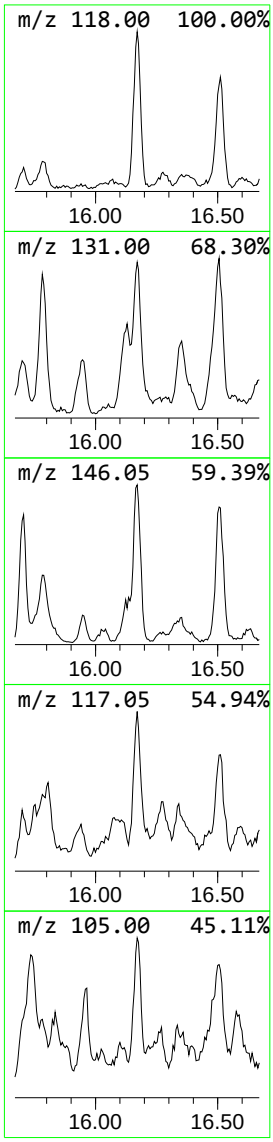
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.170	33.44 ug/l	1838650	1,4-Dichlorobenzene-d4	13.794

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	55
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	55
3	Benzene, (1-methylenebutyl)-	146	C11H14	005676-32-4	53
4	5H-Cyclohepta[b]pyridine, 6,7,8,...	163	C10H13NO	041043-09-8	46
5	2,3,4,5,6,7-Hexahydro-1H-cyclope...	146	C11H14	1010189-31-0	45





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN021424\
Data File : VN081032.D
Acq On : 14 Feb 2024 18:07
Operator : JC\MD
Sample : P1459-03 20X
Misc : 1.01g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WAL-OILY-DEBRIS

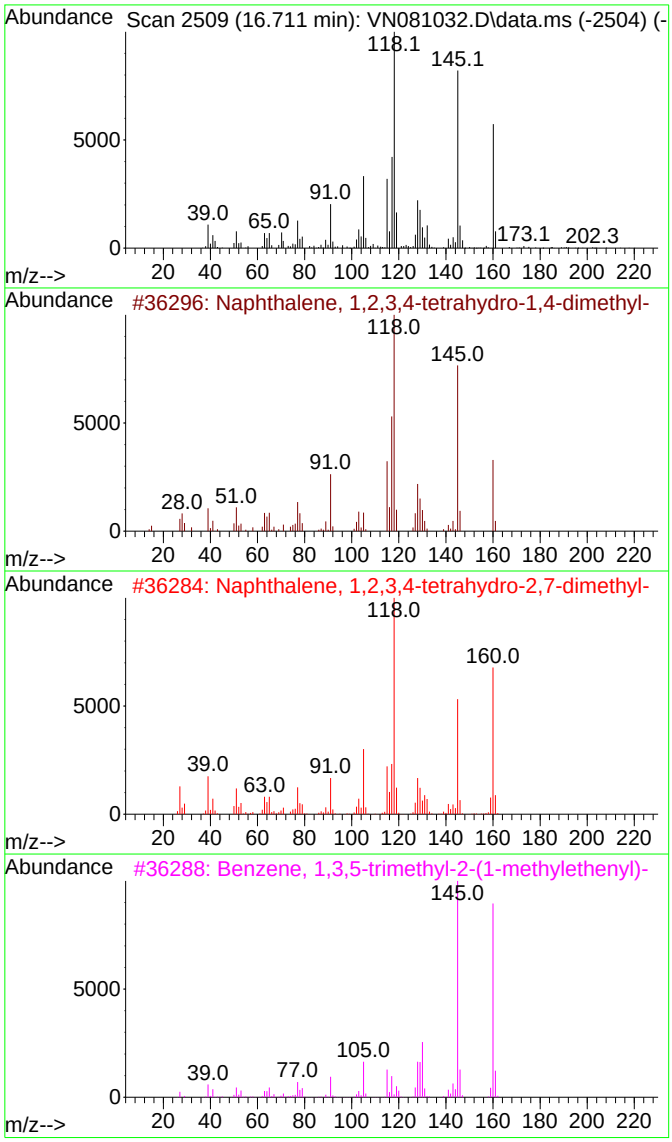
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021424W.M
Quant Title : SW846 8260

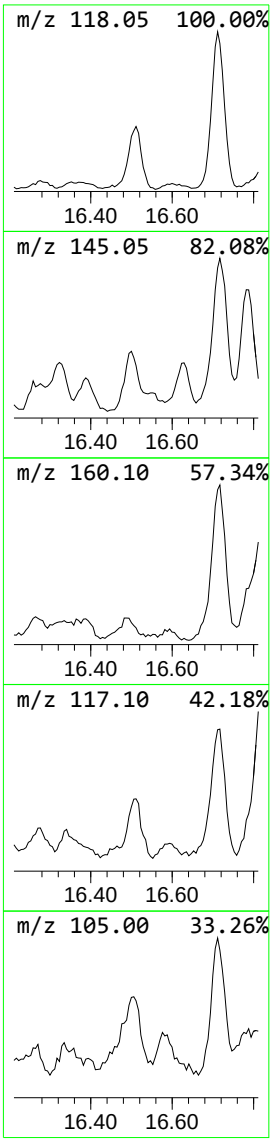
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.711	32.05 ug/l	1762340	1,4-Dichlorobenzene-d4	13.794

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	95
2	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	76
3	Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	70
4	Benzene, 1,2,4-trimethyl-5-(1-me...	160	C12H16	054340-84-0	70
5	1H-Indene, 2,3-dihydro-1,1,3-tri...	160	C12H16	002613-76-5	60





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN021424\
Data File : VN081032.D
Acq On : 14 Feb 2024 18:07
Operator : JC\MD
Sample : P1459-03 20X
Misc : 1.01g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WAL-OILY-DEBRIS

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021424W.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
Naphthalene, de...	14.123	24.0	ug/l	1320710	4	13.794	2749090	50.0
Naphthalene, de...	14.629	21.3	ug/l	1171970	4	13.794	2749090	50.0
2-(1-Adamantyl)...	14.664	18.3	ug/l	1005170	4	13.794	2749090	50.0
1-Methyldecahyd...	14.794	21.8	ug/l	1197440	4	13.794	2749090	50.0
1-Phenyl-1-butene	15.059	26.8	ug/l	1473070	4	13.794	2749090	50.0
unknown15.111	15.111	17.6	ug/l	968894	4	13.794	2749090	50.0
Naphthalene, 1,...	15.217	21.1	ug/l	1162980	4	13.794	2749090	50.0
Naphthalene, 1,...	15.700	39.8	ug/l	2189830	4	13.794	2749090	50.0
Naphthalene, 1,...	16.170	33.4	ug/l	1838650	4	13.794	2749090	50.0
Naphthalene, 1,...	16.711	32.1	ug/l	1762340	4	13.794	2749090	50.0