LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM020818\
Data File : BM014197.D
Acq On : 08 Feb 2018 17:34
Operator : SJ/JU
Sample : J1455-09
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : LW-04-B

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:\HPCHEM1\BNA_M\METHODS\8270-BM020718.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.081	12	16	21	rBV	124503	151657	1.59%	0.320%
2	5.163	364	370	384	rBV	1007755	1521493	15.98%	3.212%
3	6.734	631	637	650	rBV	564988	958409	10.07%	2.023%
4	7.081	689	696	704	rBV	2305583	3550592	37.30%	7.496%
5	7.540	768	774	782	rBV	484345	779091	8.18%	1.645%
6	7.857	821	828	834	rBV	2272141	3628809	38.12%	7.661%
7	7.904	834	836	844	rVB	105634	120188	1.26%	0.254%
8	8.698	964	971	989	rBV	1783419	3102211	32.59%	6.550%
9	9.716	1135	1144	1152	rBV4	65780	139694	1.47%	0.295%
10	10.316	1238	1246	1260	rBV	565203	1070469	11.24%	2.260%
11	12.810	1662	1670	1684	rBV	4182189	6401710	67.24%	13.516%
12	13.504	1782	1788	1794	rBV10	35318	103440	1.09%	0.218%
13	14.016	1870	1875	1886	rVB10	61935	158285	1.66%	0.334%
14	14.198	1900	1906	1913	rBV2	896639	1413004	14.84%	2.983%
15	14.421	1934	1944	1947	rBV7	114601	236969	2.49%	0.500%
16	14.721	1991	1995	1998	rBV3	94438	134910	1.42%	0.285%
17	14.757	1998	2001	2006	rVB2	125417	164927	1.73%	0.348%
18	14.816	2006	2011	2014	rBV3	88734	152975	1.61%	0.323%
19	14.851	2014	2017	2024	rVB5	80390	127493	1.34%	0.269%
20	15.045	2043	2050	2054	rBV5	118688	288970	3.04%	0.610%
21	15.104	2055	2060	2063	rVV4	180619	364250	3.83%	0.769%
22	15.139	2063	2066	2075	rVB3	231750	418368	4.39%	0.883%
23	15.480	2120	2124	2127	rBV	88679	150260	1.58%	0.317%
24	15.604	2137	2145	2151	rBV	818153	1709183	17.95%	3.609%
25	15.698	2155	2161	2168	rVV	3647193	5130571	53.89%	10.832%
26	16.957	2369	2375	2381	rBV2	1120433	1623611	17.05%	3.428%
27	17.862	2525	2529	2535	rBV2	218103	309057	3.25%	0.652%
28	19.156	2746	2749	2753	rBV5	119823	147186	1.55%	0.311%
29	19.603	2819	2825	2830	rBV	7664521	9520047	100.00%	20.099%
30	20.880	3038	3042	3049	rBV2	179407	239088	2.51%	0.505%
31	21.162	3085	3090	3097	rBV	1472863	1903309	19.99%	4.018%
32	23.339	3453	3460	3473	rVB	871327	1645236	17.28%	3.473%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM020818\
Data File : BM014197.D
Acq On : 08 Feb 2018 17:34
Operator : SJ/JU
Sample : J1455-09
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : LW-04-B

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:\HPCHEM1\BNA_M\METHODS\8270-BM020718.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

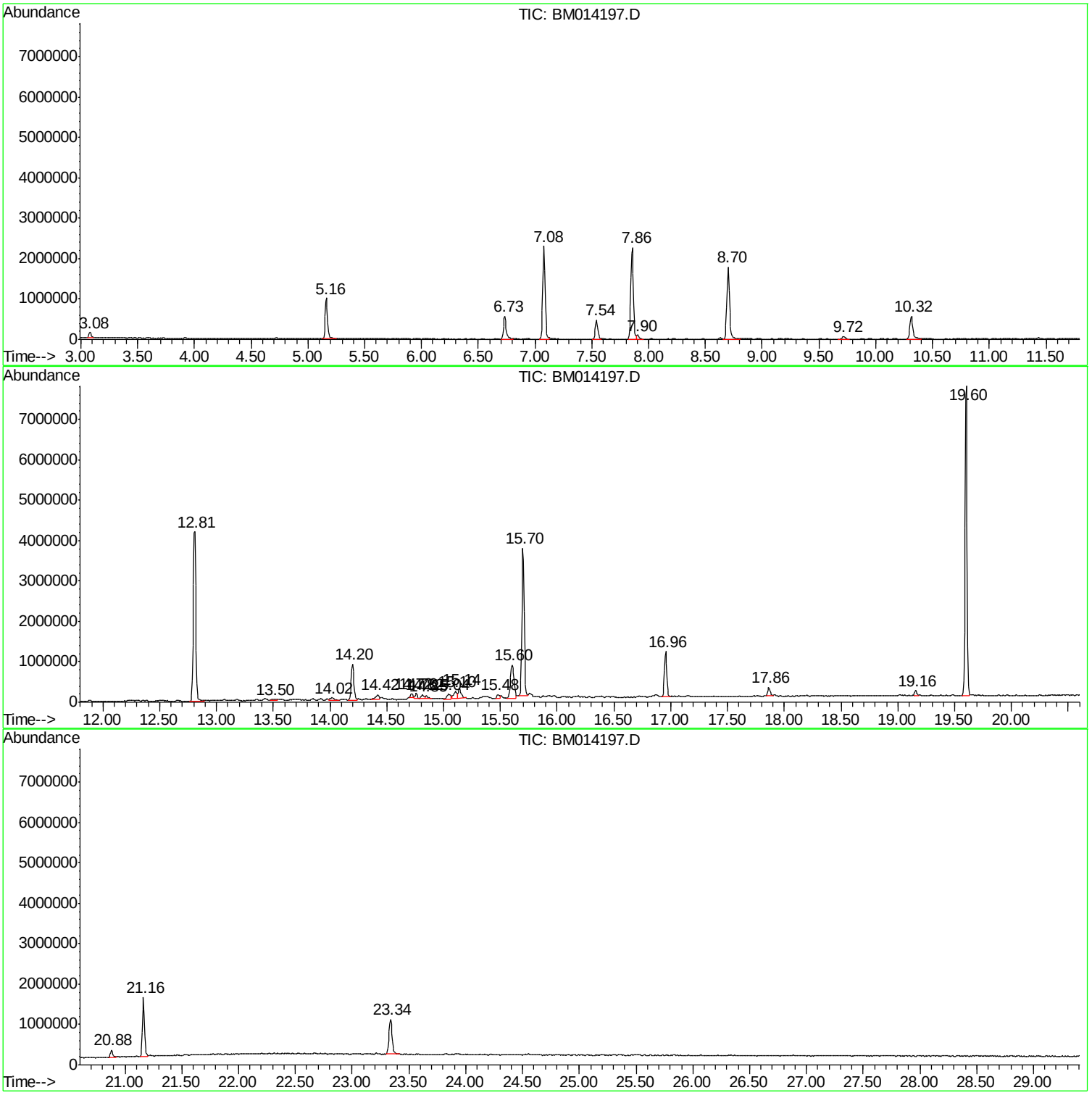
Sum of corrected areas: 47365462

Data Path : Z:\HPCHEM1\BNA M\DATA\BM020818\
Data File : BM014197.D
Acq On : 08 Feb 2018 17:34
Operator : SJ/JU
Sample : J1455-09
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : LW-04-B

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM020718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM020818\
Data File : BM014197.D
Acq On : 08 Feb 2018 17:34
Operator : SJ/JU
Sample : J1455-09
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : LW-04-B

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM020718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

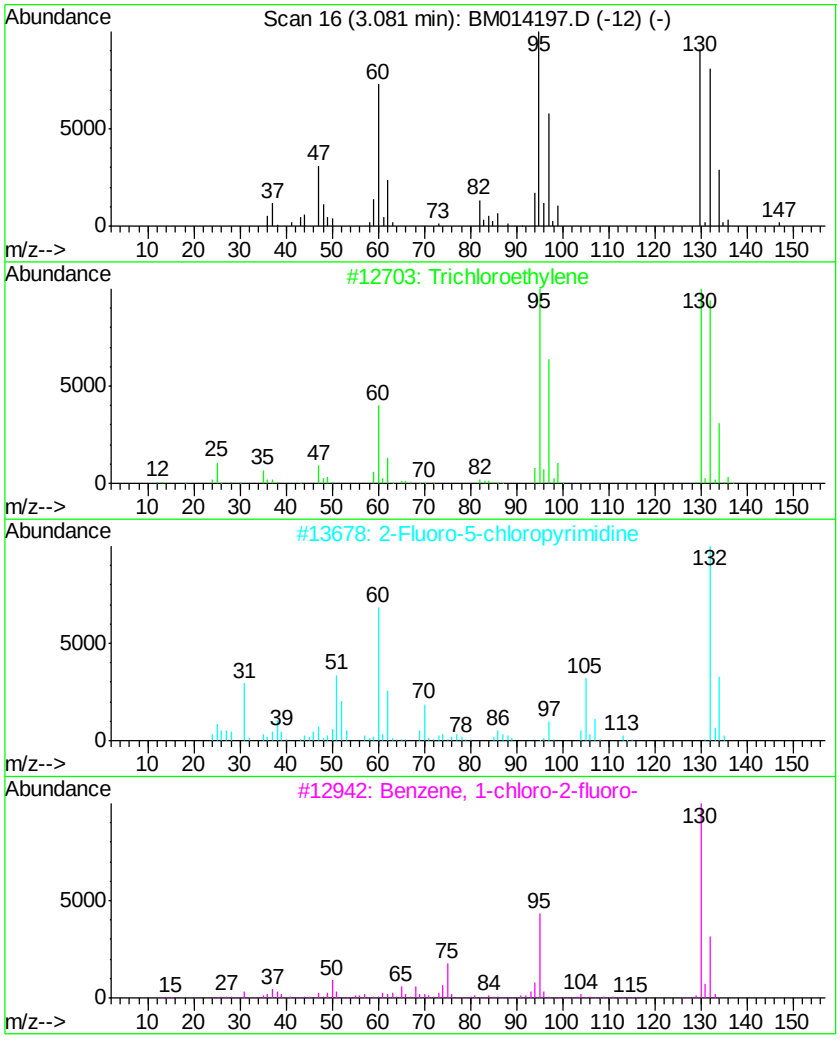
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

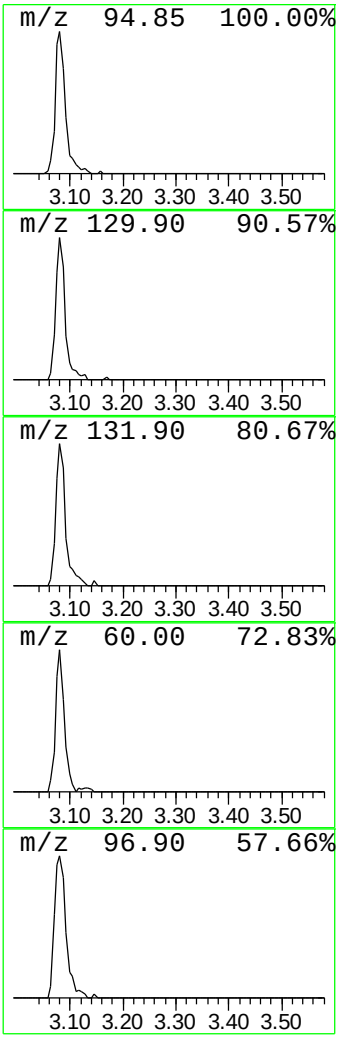
****

Peak Number 1 Trichloroethylene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.08	3.89 ng	151657	1,4-Dichlorobenzene-d4	7.54

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Trichloroethylene	130	C2HCl3	000079-01-6	93
2	2-Fluoro-5-chloropyrimidine	132	C4H2ClFN2	062802-37-3	35
3	Benzene, 1-chloro-2-fluoro-	130	C6H4ClF	000348-51-6	25
4	Thiophene, 2-(chloromethyl)-	132	C5H5ClS	000765-50-4	18
5	4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	18





Data Path : Z:\HPCHEM1\BNA M\DATA\BM020818\
Data File : BM014197.D
Acq On : 08 Feb 2018 17:34
Operator : SJ/JU
Sample : J1455-09
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : LW-04-B

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM020718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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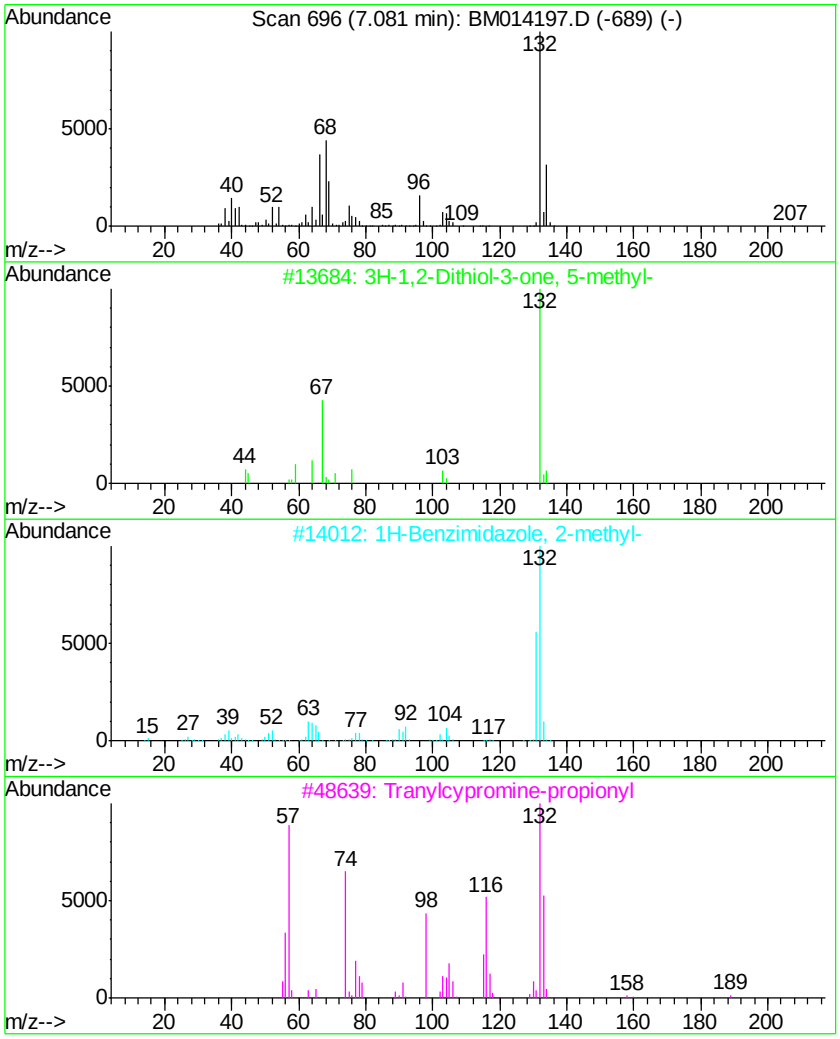
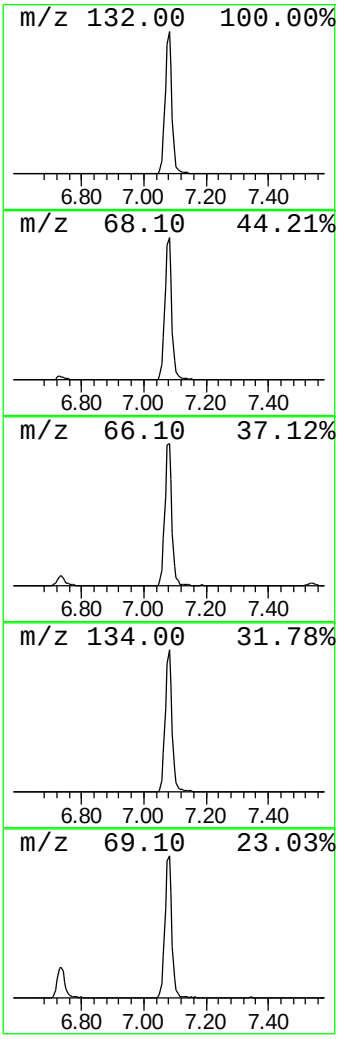
Peak Number 2 unknown7.08 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.08	91.15 ng	3550590	1,4-Dichlorobenzene-d4	7.54

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3H-1,2-Dithiol-3-one, 5-methyl-	132	C4H4OS2	003620-08-4	27
2	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	27
3	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	22
4	2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	22
5	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	17





Data Path : Z:\HPCHEM1\BNA M\DATA\BM020818\
Data File : BM014197.D
Acq On : 08 Feb 2018 17:34
Operator : SJ/JU
Sample : J1455-09
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : LW-04-B

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM020718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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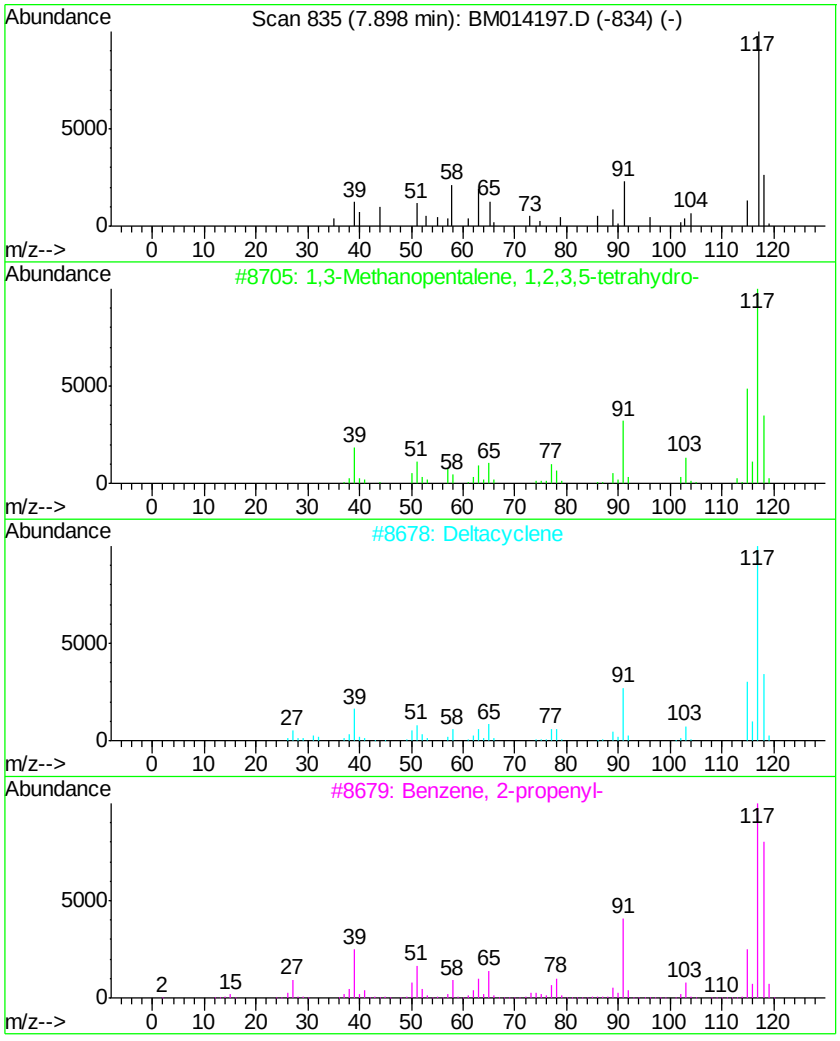
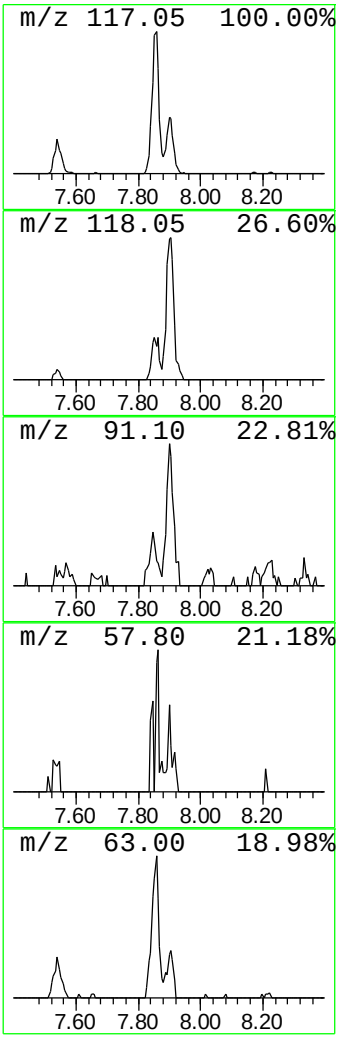
Peak Number 4 1,3-Methanopentalene, 1,2,3... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.90	3.09 ng	120188	1,4-Dichlorobenzene-d4	7.54

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Methanopentalene, 1,2,3,5-te...	118	C9H10	128600-88-4	53
2	Deltacyclene	118	C9H10	007785-10-6	50
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	45
4	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	42
5	Indane	118	C9H10	000496-11-7	42





Data Path : Z:\HPCHEM1\BNA M\DATA\BM020818\
Data File : BM014197.D
Acq On : 08 Feb 2018 17:34
Operator : SJ/JU
Sample : J1455-09
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : LW-04-B

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM020718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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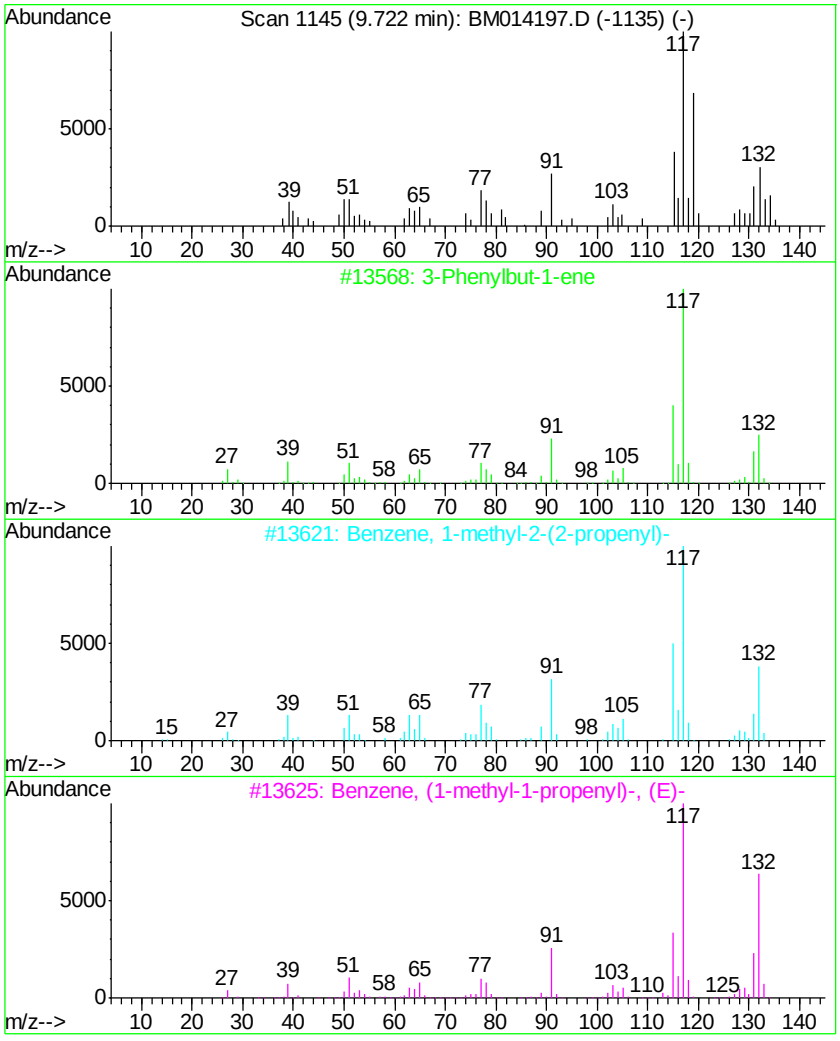
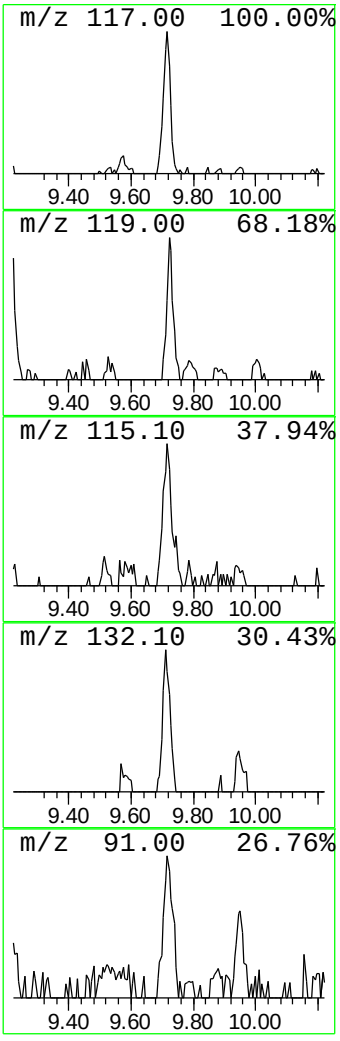
Peak Number 5 3-Phenylbut-1-ene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.72	2.61 ng	139694	Naphthalene-d8	10.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Phenylbut-1-ene	132	C10H12	000934-10-1	68
2	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	68
3	Benzene, (1-methyl-1-propenyl)-, ...	132	C10H12	000768-00-3	68
4	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	64
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	64





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Operator : SJ/JU
Sample : J1455-09
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : LW-04-B

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM020718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

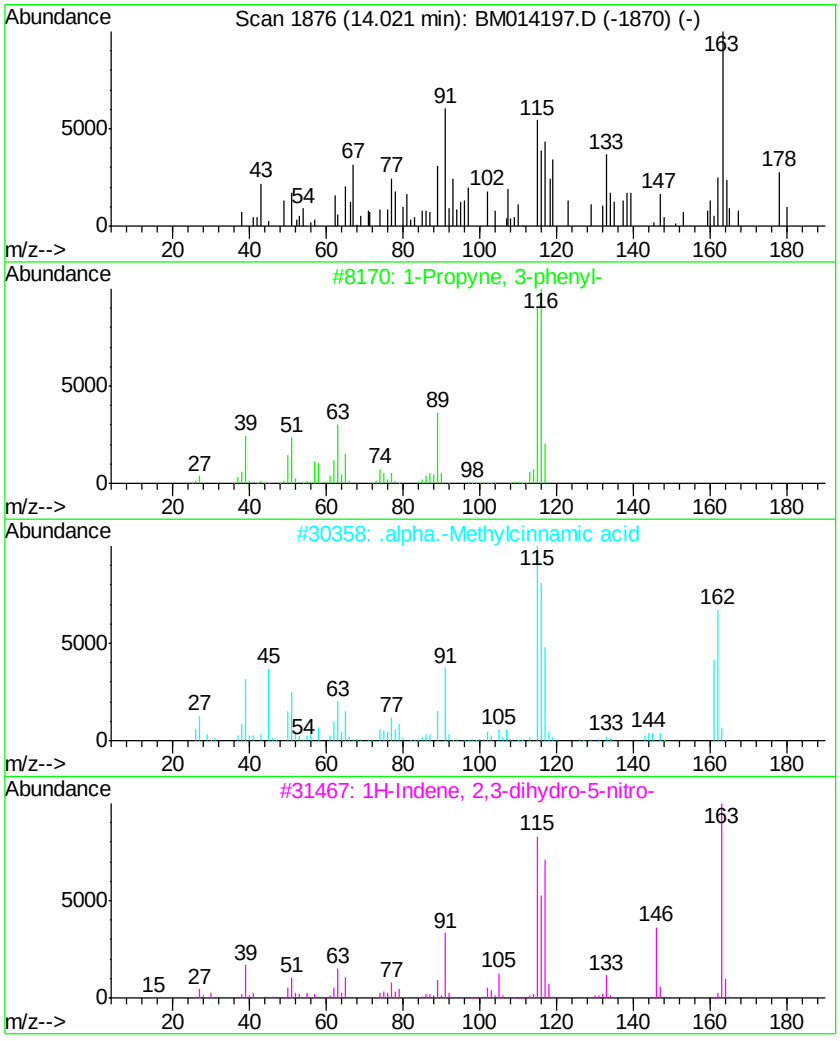
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

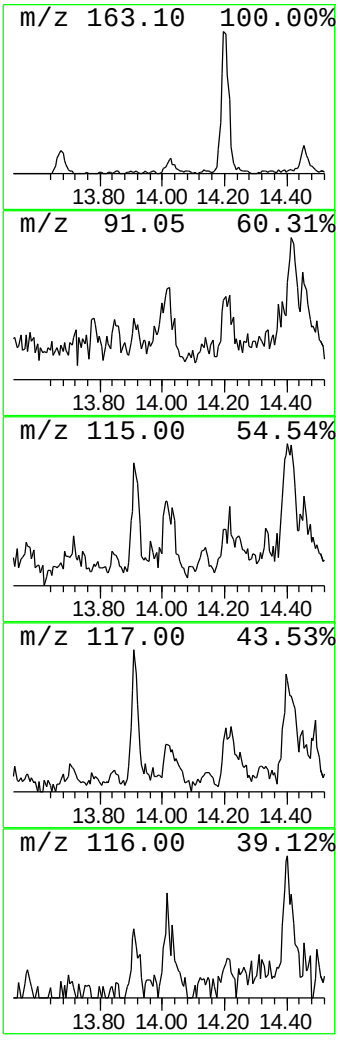
****

Peak Number 6 unknown14.02 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.02	2.24 ng	158285	Acenaphthene-d10	14.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	46
2	.alpha.-Methylcinnamic acid	162	C10H10O2	001199-77-5	35
3	1H-Indene, 2,3-dihydro-5-nitro-	163	C9H9NO2	007436-07-9	27
4	Benzene, 1-cyclopropyl-4-nitro-	163	C9H9NO2	006921-44-4	27
5	Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	27





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Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : LW-04-B

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM020718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

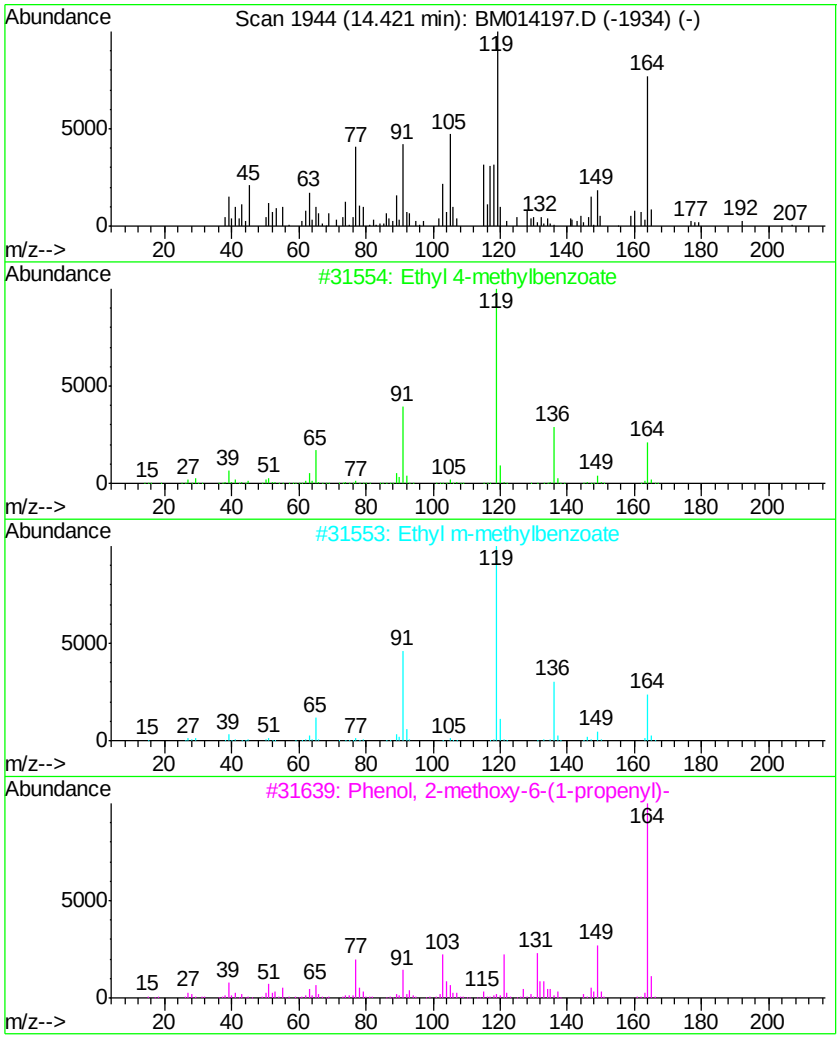
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

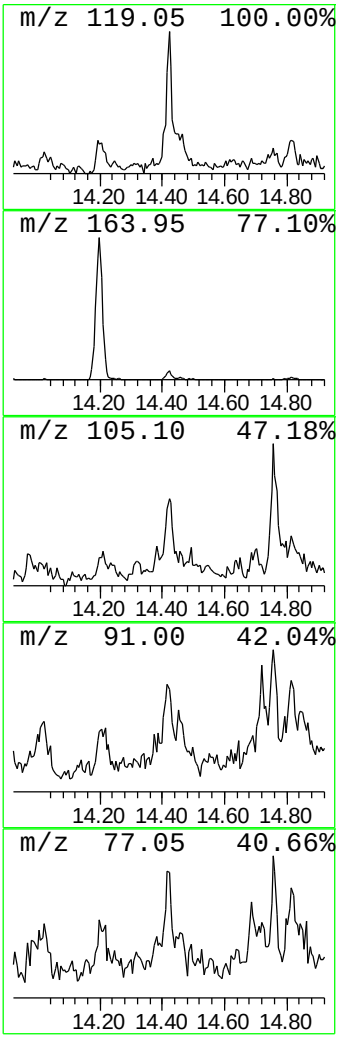
****

Peak Number 7 unknown14.42 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.42	3.35 ng	236969	Acenaphthene-d10	14.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethyl 4-methylbenzoate	164	C10H12O2	000094-08-6	49
2	Ethyl m-methylbenzoate	164	C10H12O2	000120-33-2	46
3	Phenol, 2-methoxy-6-(1-propenyl)-	164	C10H12O2	001076-55-7	43
4	Phenol, 2-methoxy-4-(1-propenyl)-	164	C10H12O2	000097-54-1	38
5	Pyridine, 5-ethenyl-2-methyl-	119	C8H9N	000140-76-1	38





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ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : LW-04-B

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM020718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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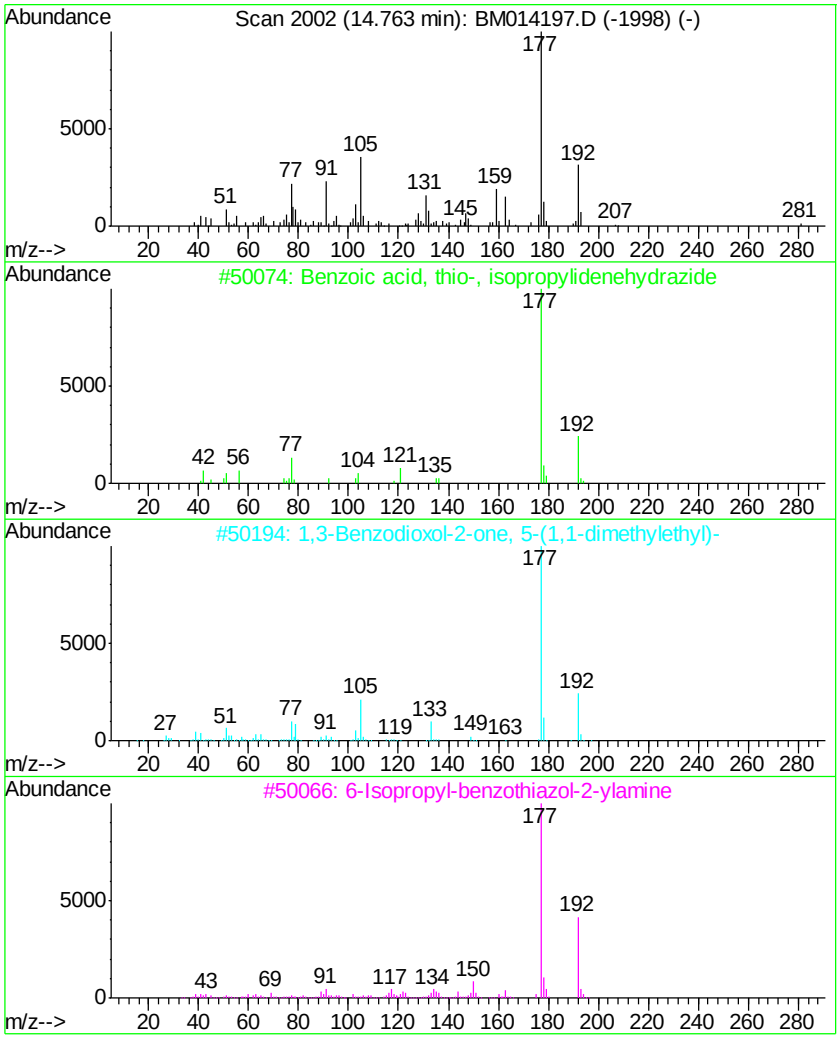
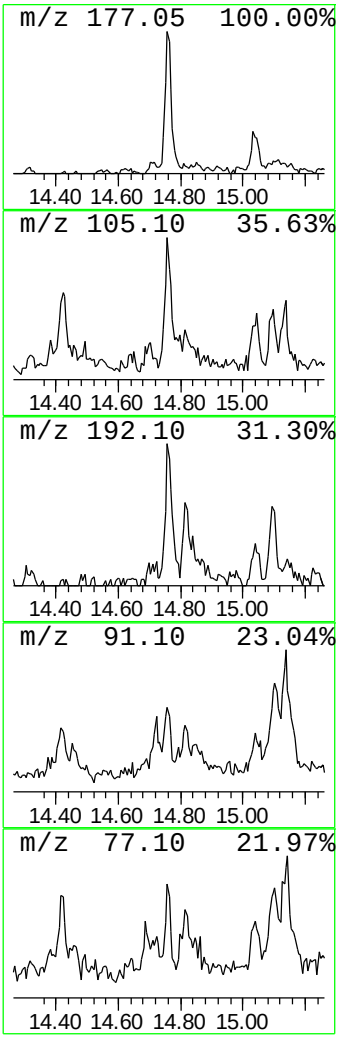
Peak Number 8 Benzoic acid, thio-, isopro... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.76	2.33 ng	164927	Acenaphthene-d10	14.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, thio-, isopropylid...	192	C10H12N2S	020185-02-8	52
2	1,3-Benzodioxol-2-one, 5-(1,1-di...	192	C11H12O3	054815-21-3	52
3	6-Isopropyl-benzothiazol-2-ylamine	192	C10H12N2S	032895-14-0	49
4	5-(4,5-Dihydro-3H-pyrrol-2-ylmet...	192	C11H16N2O	051430-87-6	47
5	Pyrazine, 2,5-bis(1,1-dimethylet...	192	C12H20N2	018709-51-8	47





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ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM020718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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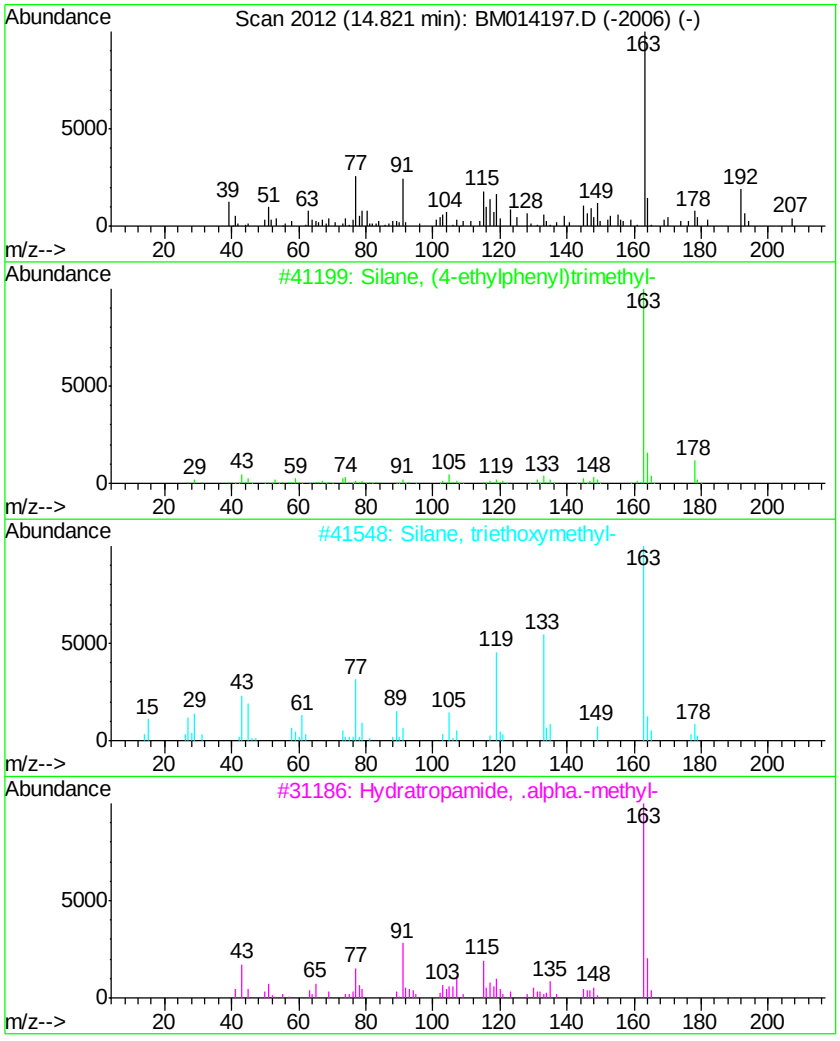
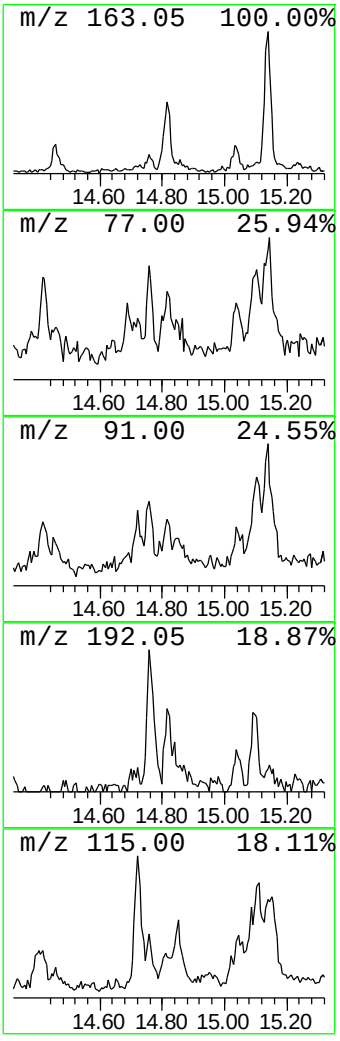
Peak Number 9 Silane, (4-ethylphenyl)trim... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	2.17 ng	152975	Acenaphthene-d10	14.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Silane, (4-ethylphenyl)trimethyl-	178	C11H18Si	017988-50-0	58
2	Silane, triethoxymethyl-	178	C7H18O3Si	002031-67-6	53
3	Hydratropamide, .alpha.-methyl-	163	C10H13NO	000826-54-0	52
4	Benzoic acid, p-tert-butyl-	178	C11H14O2	000098-73-7	52
5	Propofol	178	C12H18O	002078-54-8	50





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Sample : J1455-09
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : LW-04-B

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM020718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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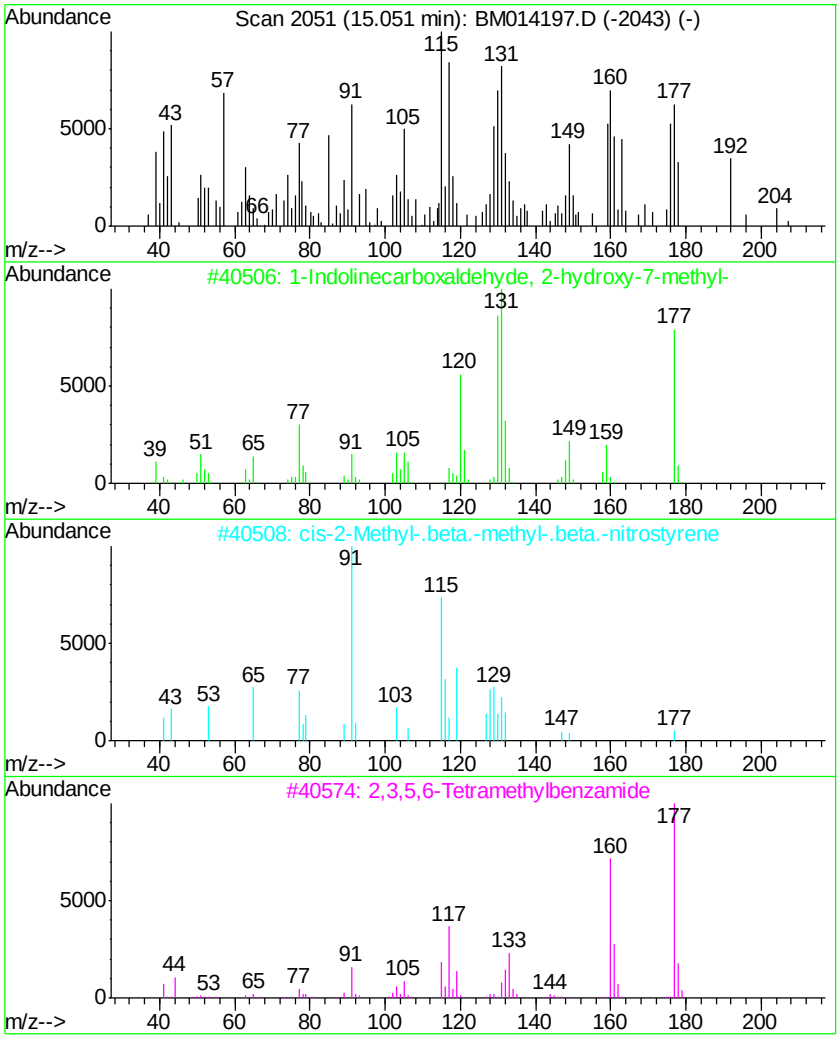
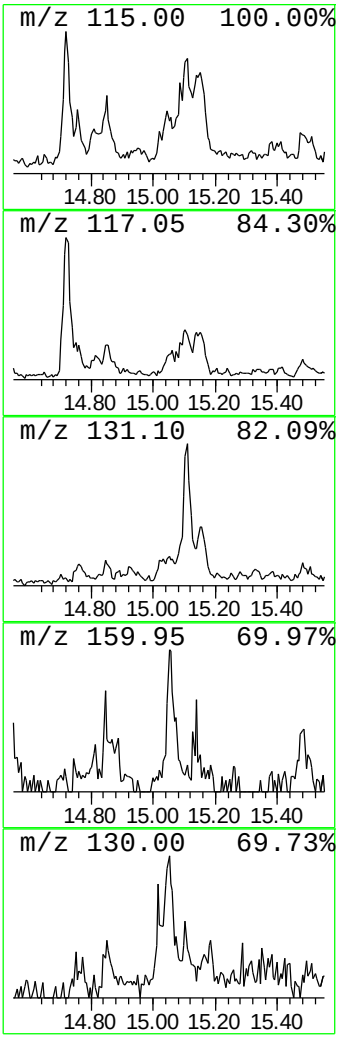
Peak Number 10 unknown15.05 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.05	4.09 ng	288970	Acenaphthene-d10	14.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Indolinecarboxaldehyde, 2-hydr...	177	C10H11NO2	022614-65-9	22
2	cis-2-Methyl-.beta.-methyl-.beta...	177	C10H11NO2	1000122-79-5	22
3	2,3,5,6-Tetramethylbenzamide	177	C11H15NO	099858-56-7	14
4	2-(1-Cyclohexenyl)cyclohexanone	178	C12H18O	001502-22-3	11
5	Pyrrole, 1-methyl-2-(trifluorome...	149	C6H6F3N	067095-61-8	11





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Data File : BM014197.D
Acq On : 08 Feb 2018 17:34
Operator : SJ/JU
Sample : J1455-09
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : LW-04-B

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM020718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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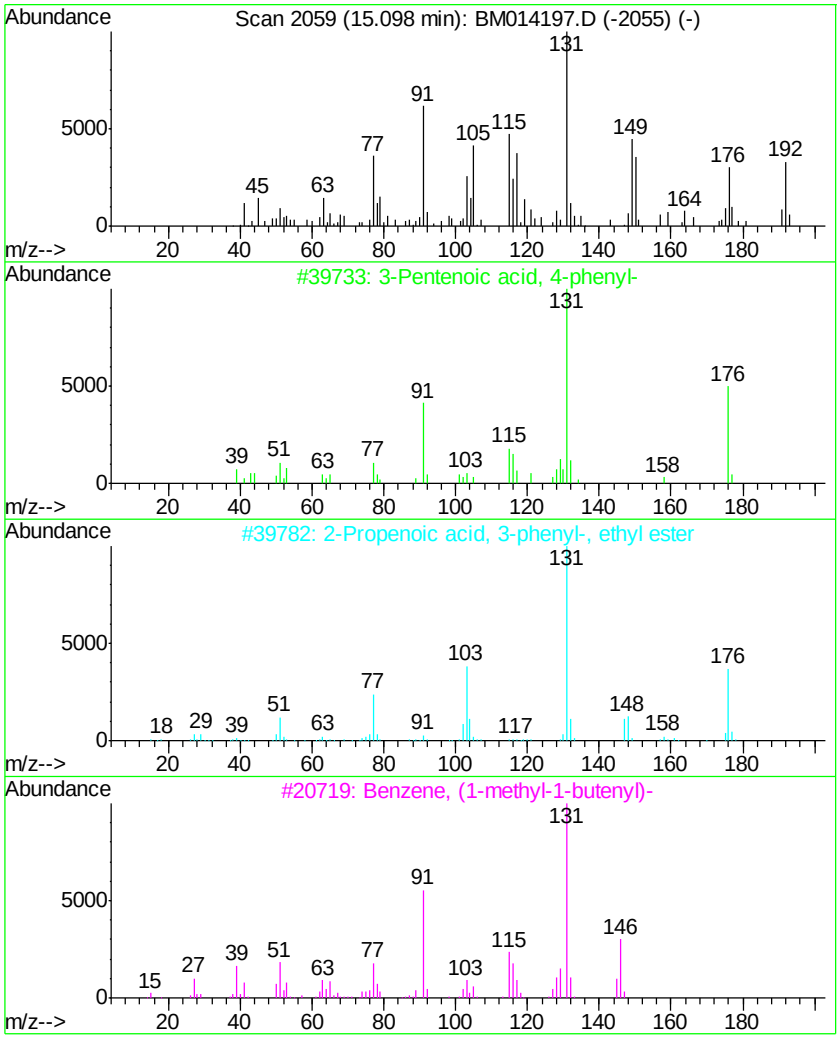
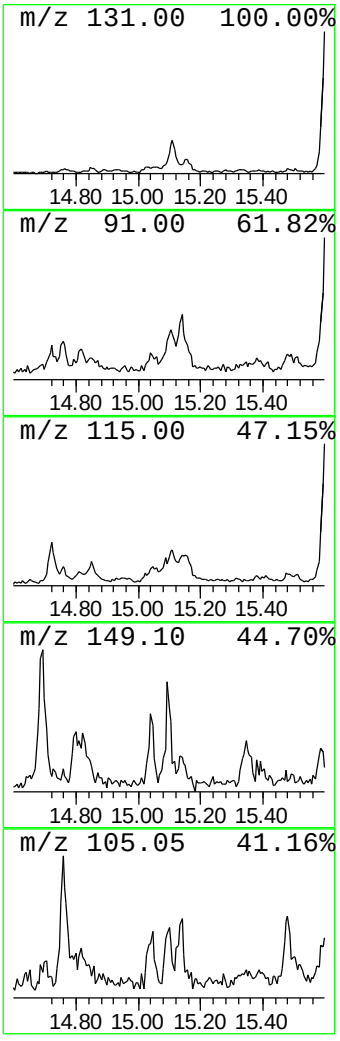
Peak Number 11 unknown15.10 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.10	5.16 ng	364250	Acenaphthene-d10	14.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Pentenoic acid, 4-phenyl-	176	C11H12O2	053774-19-9	49
2	2-Propenoic acid, 3-phenyl-, eth...	176	C11H12O2	000103-36-6	43
3	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	38
4	2-Butene, 3-chloro-1-phenyl-, (Z)-	166	C10H11Cl	016608-68-7	35
5	Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	35





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Operator : SJ/JU
Sample : J1455-09
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : LW-04-B

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM020718.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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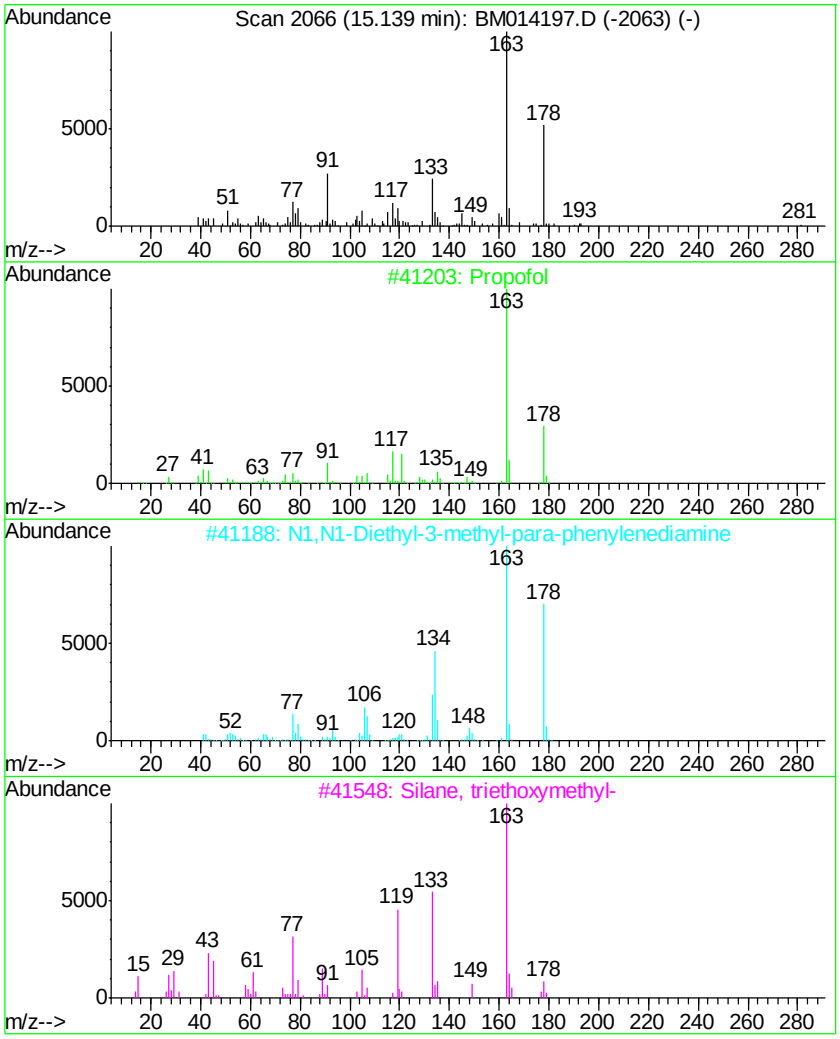
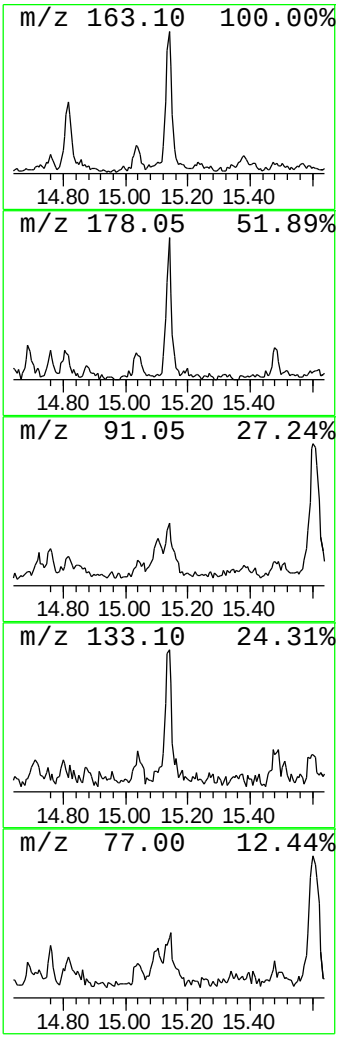
Peak Number 12 Propofol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.14	5.92 ng	418368	Acenaphthene-d10	14.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propofol	178	C12H18O	002078-54-8	64
2	N1,N1-Diethyl-3-methyl-para-phen...	178	C11H18N2	002051-79-8	59
3	Silane, triethoxymethyl-	178	C7H18O3Si	002031-67-6	53
4	Ethanone, 1-(2,3-dihydro-1,4-ben...	178	C10H10O3	002879-20-1	53
5	6-tert-Butyl-8-[(2,4-dimethoxy-b...	356	C17H20N6O3	1000295-27-2	43





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Operator : SJ/JU
Sample : J1455-09
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : LW-04-B

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-BM020718.M
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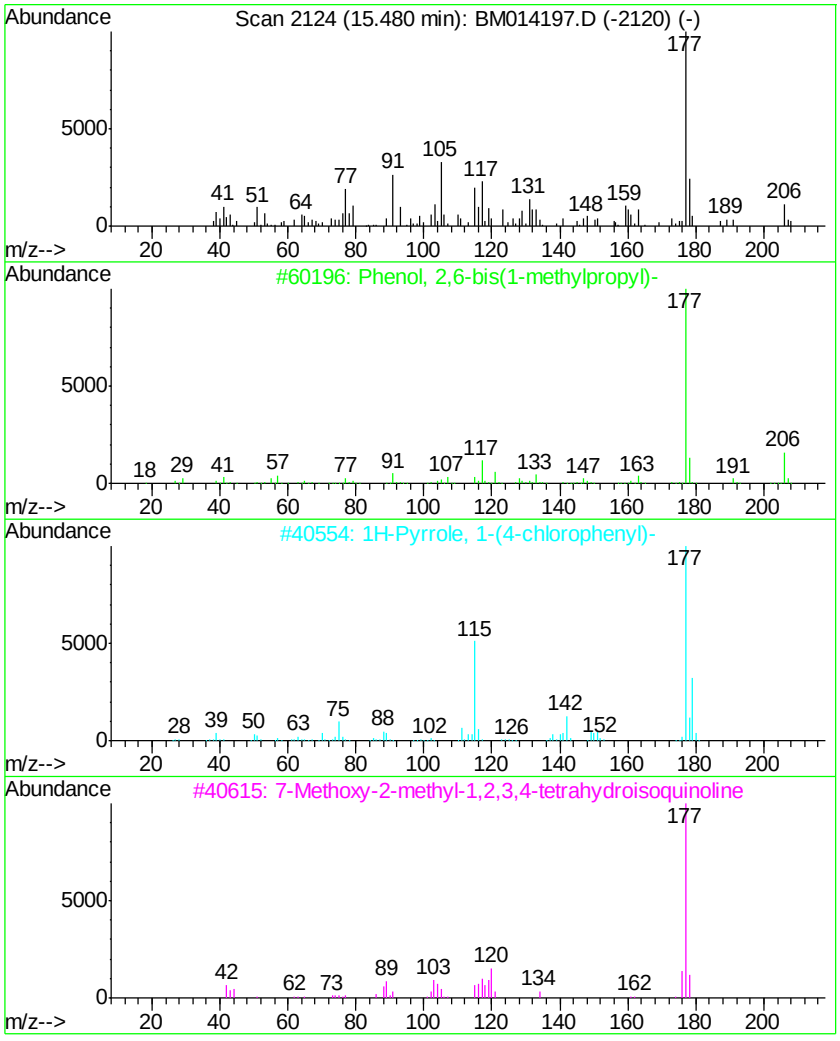
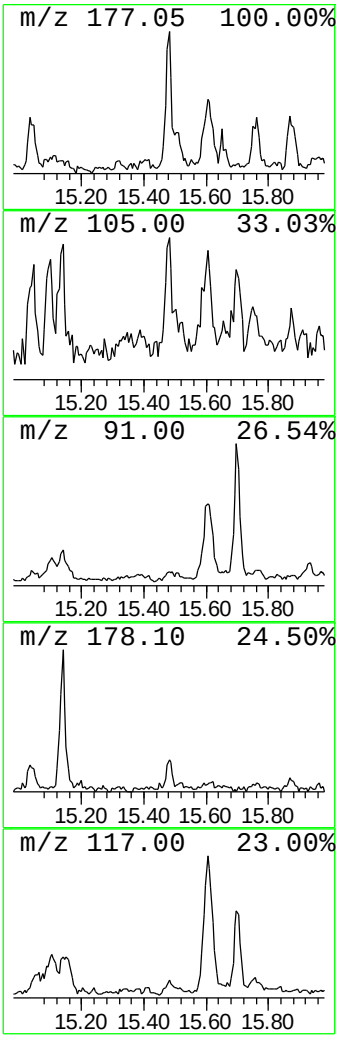
Peak Number 13 Phenol, 2,6-bis(1-methylpro... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.48	2.13 ng	150260	Acenaphthene-d10	14.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 2,6-bis(1-methylpropyl)-	206	C14H22O	005510-99-6	58
2	1H-Pyrrole, 1-(4-chlorophenyl)-	177	C10H8ClN	005044-38-2	43
3	7-Methoxy-2-methyl-1,2,3,4-tetra...	177	C11H15NO	054893-23-1	40
4	2-Ethyl-2-(p-tolyl)malonamide	220	C12H16N2O2	068692-83-1	40
5	Spiro[cyclopenta[d]-1,3,2-dioxab...	206	C12H19BO2	062238-26-0	38





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ALS Vial : 13 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : LW-04-B

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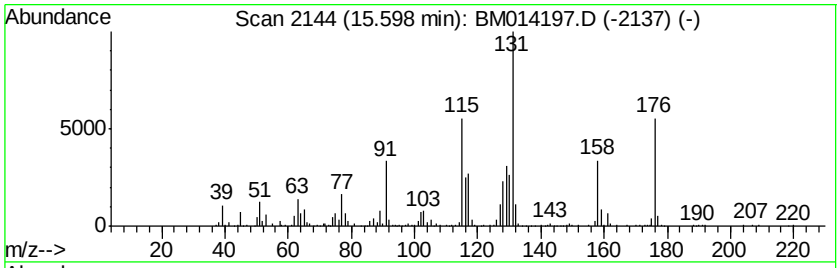
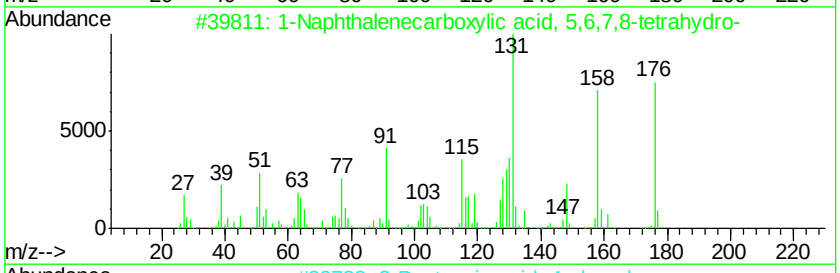
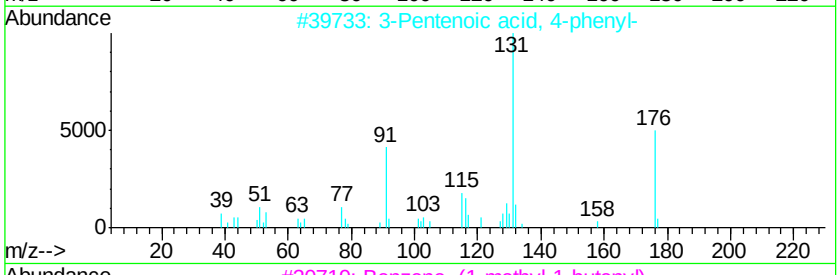
Peak Number 14 1-Naphthalenecarboxylic aci... Concentration Rank 3

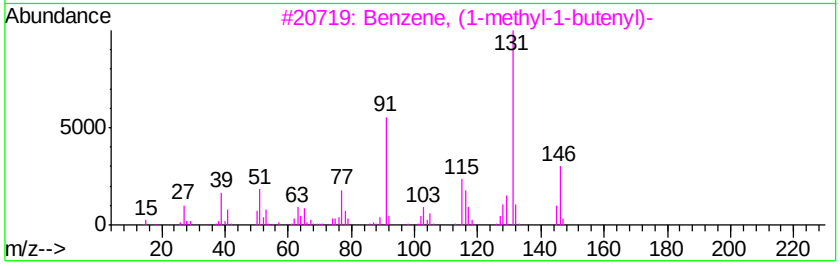
R.T.	EstConc	Area	Relative to ISTD	R.T.
15.60	21.05 ng	1709180	Phenanthrene-d10	16.96

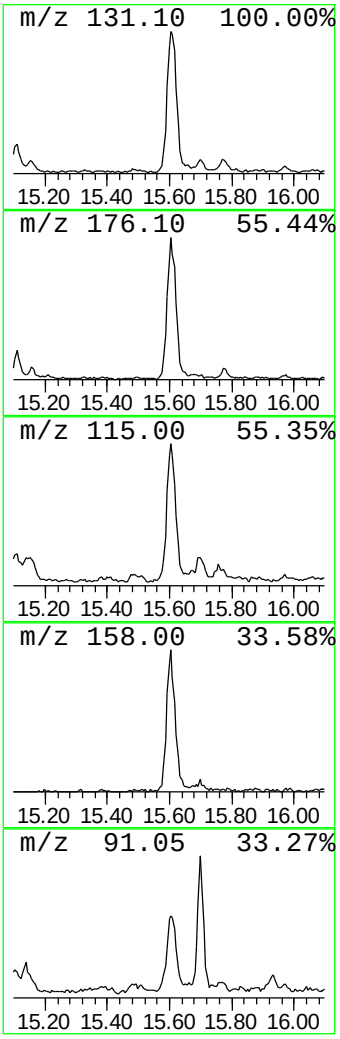
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthalenecarboxylic acid, 5,...	176	C11H12O2	004242-18-6	58
2	3-Pentenoic acid, 4-phenyl-	176	C11H12O2	053774-19-9	47
3	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	43
4	Cinnamic acid, .alpha.,.beta.-di...	176	C11H12O2	1000151-56-4	43
5	Benzene, (2-chloro-2-butenyl)-	166	C10H11Cl	054411-12-0	35











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Instrument : BNA_M
ClientSampleId : LW-04-B

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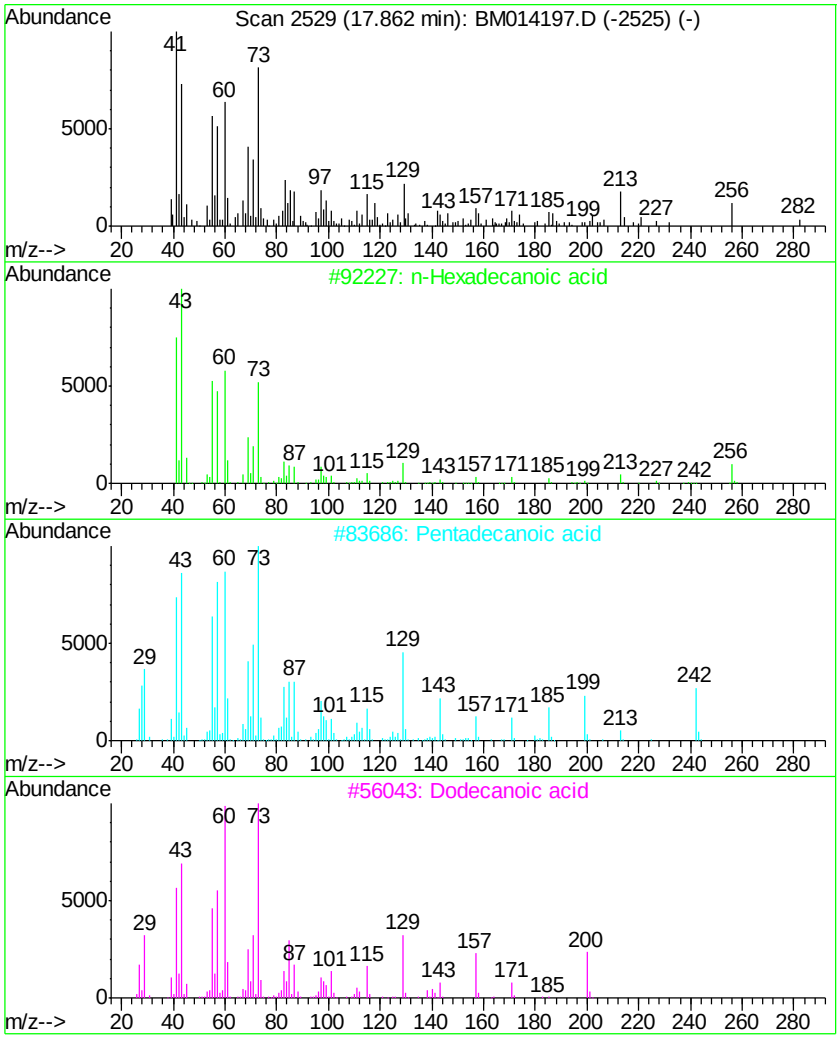
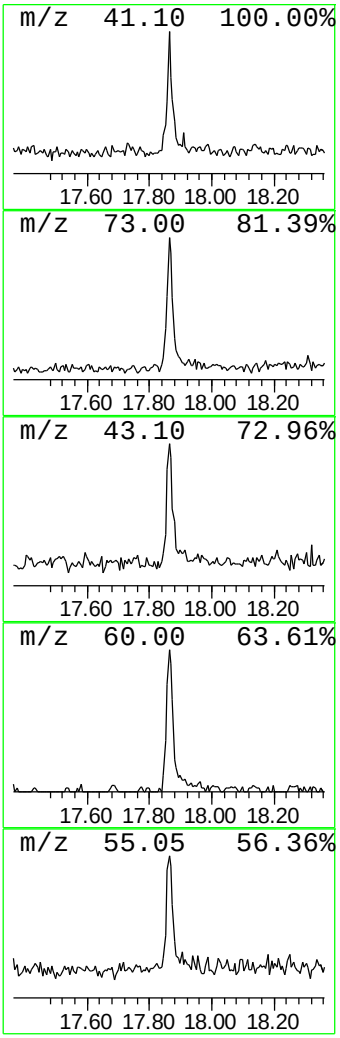
Peak Number 15 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.86	3.81 ng	309057	Phenanthrene-d10	16.96

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	58
3	Dodecanoic acid	200	C12H24O2	000143-07-7	50
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	46
5	12-Bromododecanoic acid	278	C12H23BrO2	073367-80-3	43





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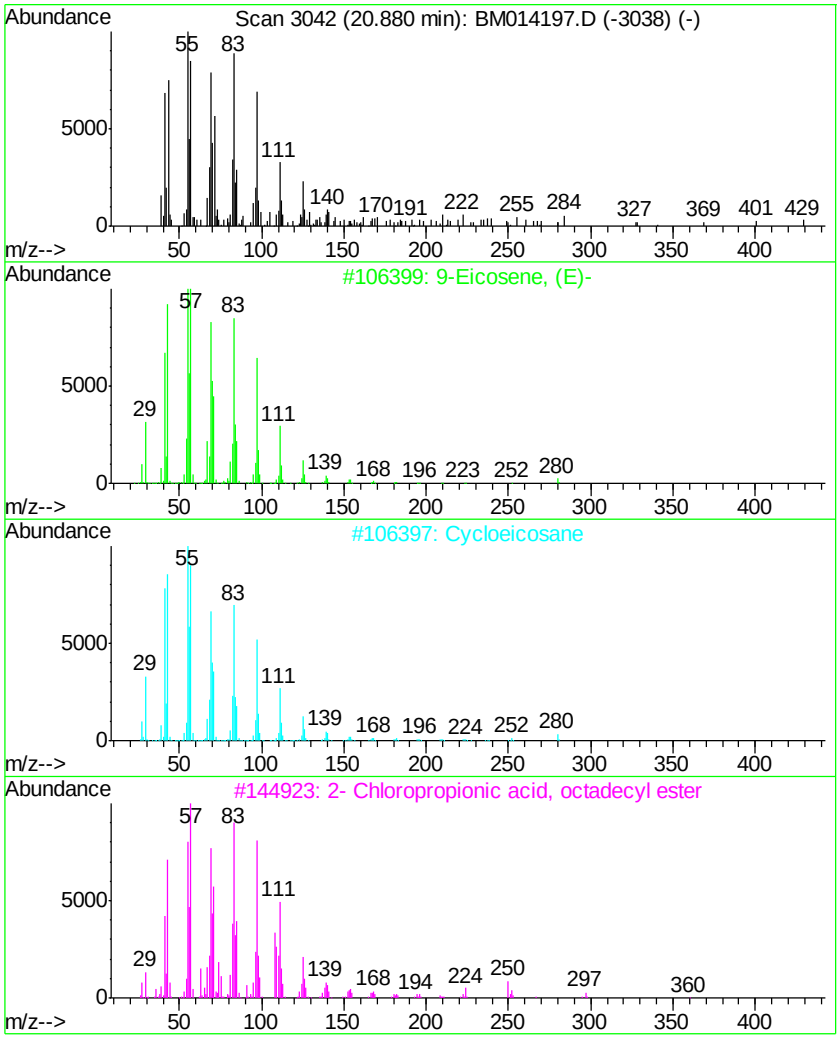
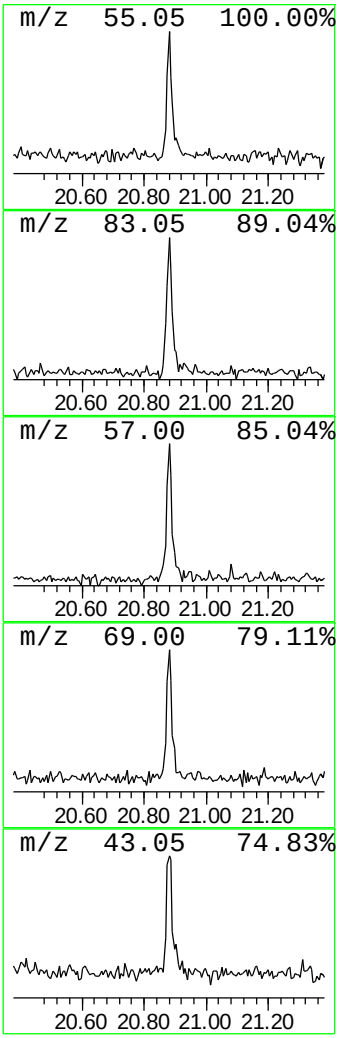
Peak Number 16 9-Eicosene, (E)- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.88	2.51 ng	239088	Chrysene-d12	21.16

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Eicosene, (E)-	280	C20H40	074685-29-3	93
2	Cycloeicosane	280	C20H40	000296-56-0	91
3	2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	91
4	17-Pentatriacontene	491	C35H70	006971-40-0	87
5	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	87





Tentatively Identified Compound (LSC) summary

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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Trichloroethylene	3.08	3.9	ng	151657	1	7.54	779091	20.0
unknown7.08	7.08	91.2	ng	3550590	1	7.54	779091	20.0
1,3-Methanopental...	7.90	3.1	ng	120188	1	7.54	779091	20.0
3-Phenylbut-1-ene	9.72	2.6	ng	139694	2	10.32	1070470	20.0
unknown14.02	14.02	2.2	ng	158285	3	14.20	1413000	20.0
unknown14.42	14.42	3.4	ng	236969	3	14.20	1413000	20.0
Benzoic acid, thi...	14.76	2.3	ng	164927	3	14.20	1413000	20.0
Silane, (4-ethylp...	14.82	2.2	ng	152975	3	14.20	1413000	20.0
unknown15.05	15.05	4.1	ng	288970	3	14.20	1413000	20.0
unknown15.10	15.10	5.2	ng	364250	3	14.20	1413000	20.0
Propofol	15.14	5.9	ng	418368	3	14.20	1413000	20.0
Phenol, 2,6-bis(1...	15.48	2.1	ng	150260	3	14.20	1413000	20.0
1-Naphthalenecarb...	15.60	21.1	ng	1709180	4	16.96	1623610	20.0
n-Hexadecanoic acid	17.86	3.8	ng	309057	4	16.96	1623610	20.0
9-Eicosene, (E)-	20.88	2.5	ng	239088	5	21.16	1903310	20.0