LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122116\
Data File : BF091867.D
Acq On : 22 Dec 2016 6:39
Operator : UM/SJ
Sample : H6156-13
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

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_F\METHODS\8270-BF122116.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	3.721	130	143	145	rBV	14163210	79757737	65.92%	12.335%
2	4.933	245	249	252	rBV	925419	1403892	1.16%	0.217%
3	5.207	269	273	275	rBV	10171550	15785260	13.05%	2.441%
4	5.356	279	286	289	rBV2	14578053	42586751	35.20%	6.586%
5	5.619	302	309	311	rVV2	15318434	56217405	46.46%	8.694%
6	6.133	351	354	356	rBV	1034362	1501499	1.24%	0.232%
7	6.201	358	360	363	rBV	2384427	2433522	2.01%	0.376%
8	6.270	363	366	367	rVV	1850096	2035849	1.68%	0.315%
9	6.304	367	369	371	rVV	3789374	4381978	3.62%	0.678%
10	6.350	371	373	375	rVV	5856720	7525612	6.22%	1.164%
11	6.407	375	378	379	rVV	2363964	3234847	2.67%	0.500%
12	6.430	379	380	386	rVV2	2657318	6763364	5.59%	1.046%
13	6.533	386	389	390	rVV	4662967	8555276	7.07%	1.323%
14	6.590	390	394	396	rVV2	12004240	25255263	20.87%	3.906%
15	6.636	396	398	400	rVB	3213782	4360948	3.60%	0.674%
16	6.762	406	409	411	rBV	988036	1716494	1.42%	0.265%
17	6.819	411	414	416	rBV	4326571	9019091	7.45%	1.395%
18	6.876	416	419	420	rVV2	2160296	3985649	3.29%	0.616%
19	6.933	420	424	428	rVB3	4026858	12812102	10.59%	1.981%
20	7.082	428	437	439	rBV2	13390524	66655705	55.09%	10.309%
21	7.173	442	445	448	rBV2	771314	1594799	1.32%	0.247%
22	7.333	453	459	462	rBV2	3097761	8032927	6.64%	1.242%
23	7.539	473	477	478	rBV2	658698	1471619	1.22%	0.228%
24	7.813	495	501	503	rBV3	9107595	28648598	23.68%	4.431%
25	7.927	508	511	517	rVB2	1766497	4643976	3.84%	0.718%
26	8.179	519	533	535	rVB2	14808171	120990288	100.00%	18.712%
27	8.487	557	560	563	rBV3	804465	2490454	2.06%	0.385%
28	8.545	563	565	567	rVB	1150120	1475882	1.22%	0.228%
29	8.647	571	574	575	rVB2	1497837	1977909	1.63%	0.306%
30	8.705	575	579	581	rBV2	1017928	1823701	1.51%	0.282%
31	8.773	581	585	587	rBV	11825348	25108353	20.75%	3.883%
32	8.876	589	594	596	rBV	10956309	24185203	19.99%	3.740%
33	9.116	611	615	617	rVV	8561682	9397283	7.77%	1.453%
34	9.219	620	624	626	rBV	4215079	6025190	4.98%	0.932%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122116\
Data File : BF091867.D
Acq On : 22 Dec 2016 6:39
Operator : UM/SJ
Sample : H6156-13
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

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_F\METHODS\8270-BF122116.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	9.379	634	638	640	rVV	1747908	2820852	2.33%	0.436%
36	9.448	642	644	645	rVV	3504350	3662057	3.03%	0.566%
37	9.470	645	646	648	rVV2	1686447	1789419	1.48%	0.277%
38	9.562	651	654	658	rVB	1723902	2143072	1.77%	0.331%
39	9.653	658	662	664	rVB	7343472	8574753	7.09%	1.326%
40	9.779	669	673	675	rBV2	2066907	3668549	3.03%	0.567%
41	9.813	675	676	678	rVV	2057973	1982314	1.64%	0.307%
42	10.042	693	696	698	rVB	2098525	2220326	1.84%	0.343%
43	10.579	739	743	745	rBV2	4422009	6359851	5.26%	0.984%
44	11.265	800	803	804	rVV	3010666	3314371	2.74%	0.513%
45	11.288	804	805	807	rVV	4868803	3831468	3.17%	0.593%
46	12.854	939	942	945	rBV	7125591	7919329	6.55%	1.225%
47	13.894	1030	1033	1036	rBV	2642537	2513417	2.08%	0.389%
48	15.288	1152	1155	1158	rBV	1448692	1953145	1.61%	0.302%

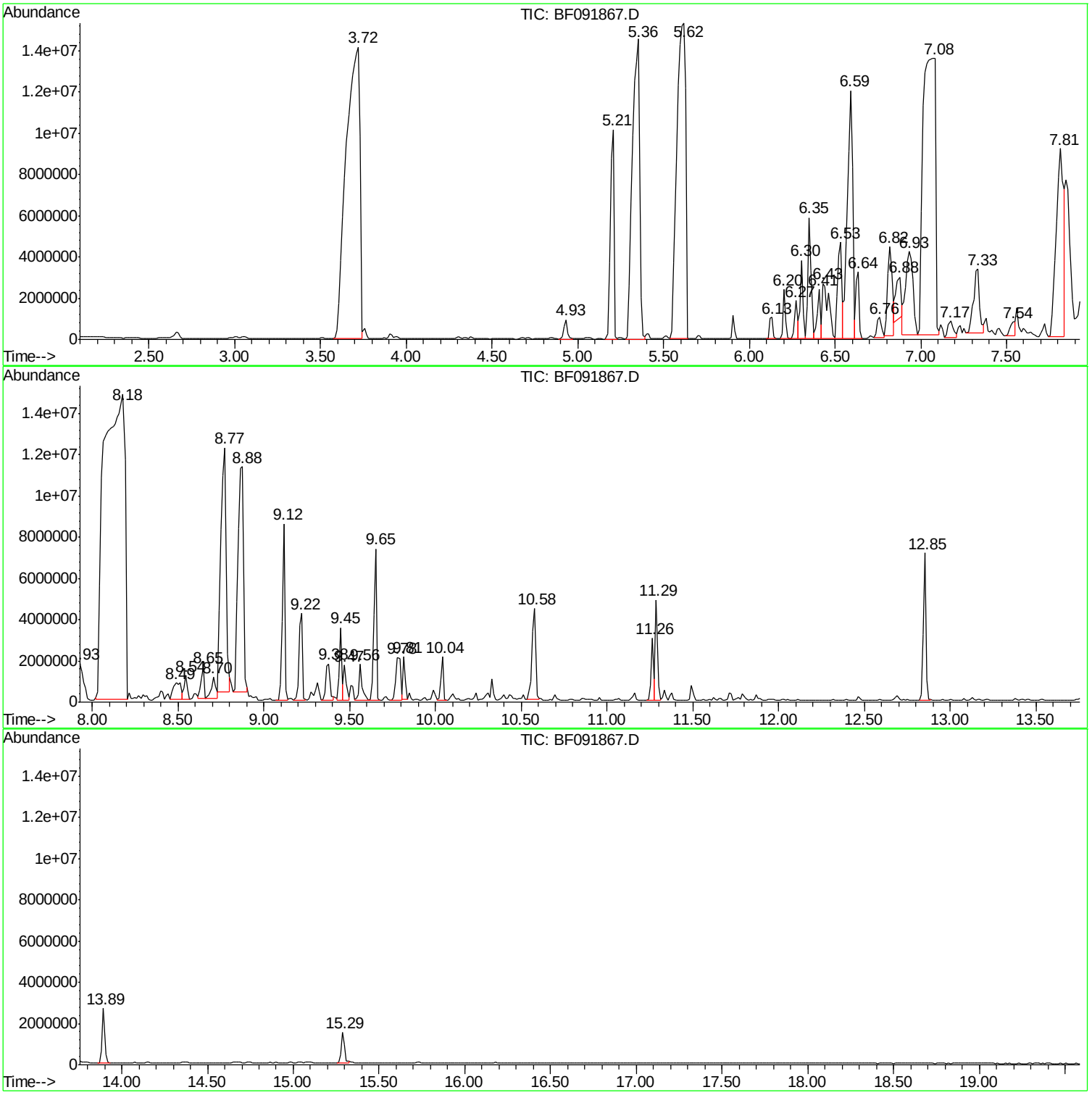
Sum of corrected areas: 646607349

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122116\
Data File : BF091867.D
Acq On : 22 Dec 2016 6:39
Operator : UM/SJ
Sample : H6156-13
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA F\DATA\BF122116\
Data File : BF091867.D
Acq On : 22 Dec 2016 6:39
Operator : UM/SJ
Sample : H6156-13
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.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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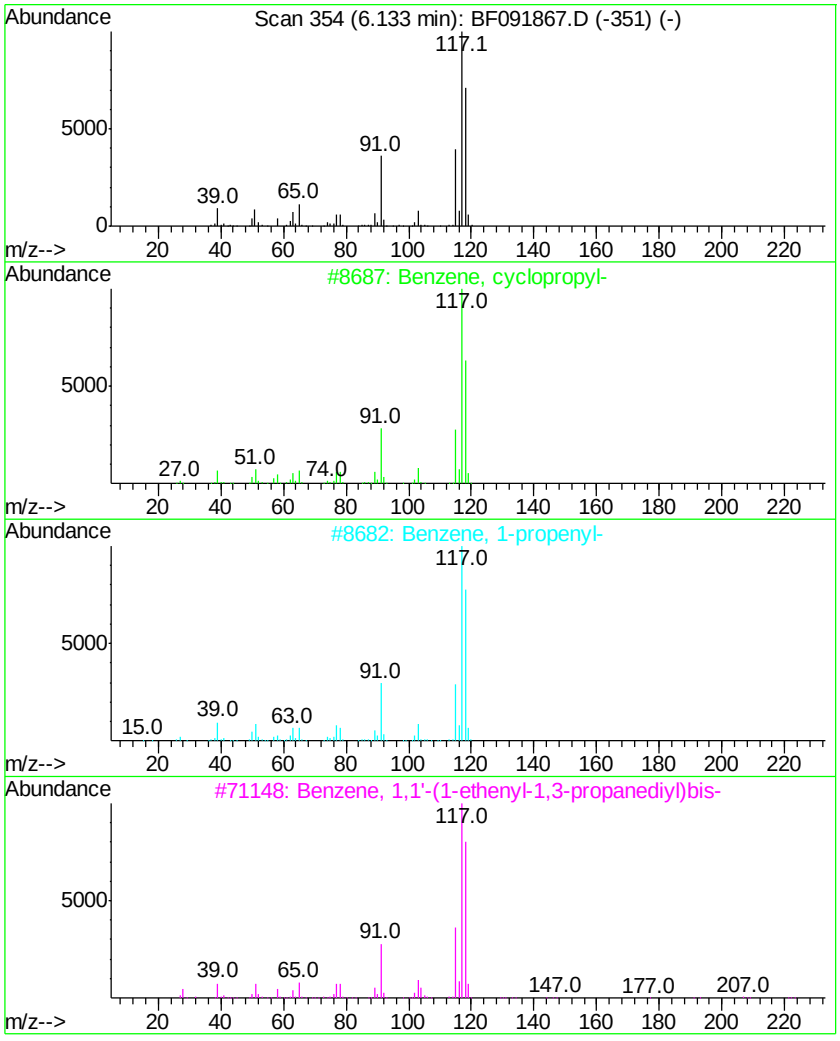
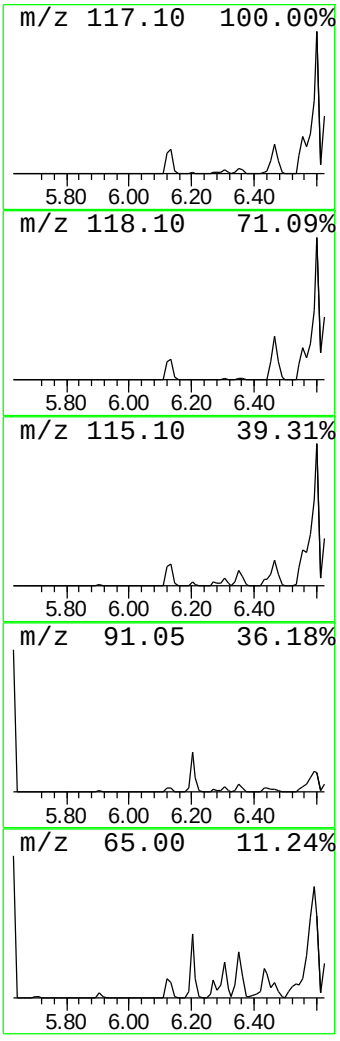
Peak Number 4 Benzene, cyclopropyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.13	17.50 ng	1501500	1,4-Dichlorobenzene-d4	6.76

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, cyclopropyl-	118	C9H10	000873-49-4	95
2	Benzene, 1-propenyl-	118	C9H10	000637-50-3	93
3	Benzene, 1,1'-(1-ethenyl-1,3-pro...	222	C17H18	061141-97-7	90
4	Benzene, 2-propenyl-	118	C9H10	000300-57-2	89
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	76





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122116\
Data File : BF091867.D
Acq On : 22 Dec 2016 6:39
Operator : UM/SJ
Sample : H6156-13
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.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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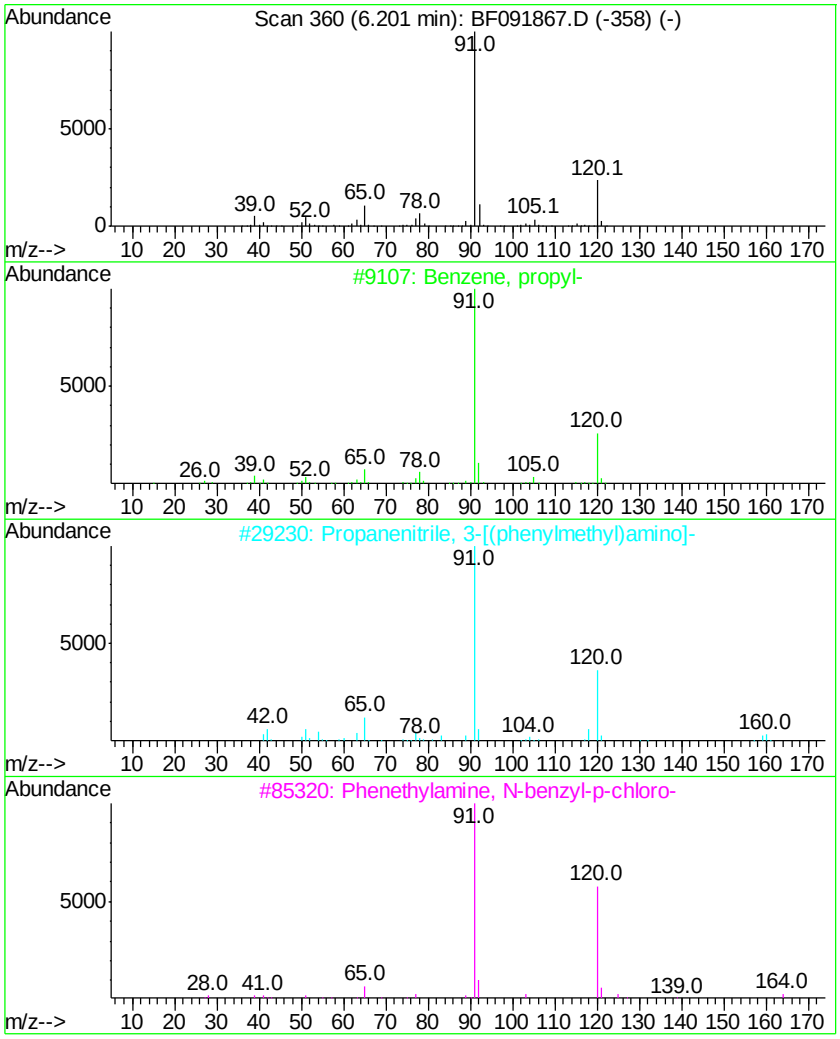
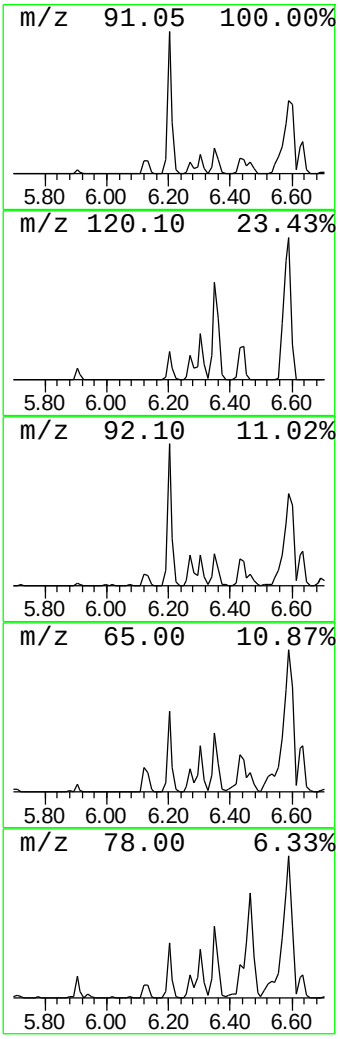
Peak Number 5 Benzene, propyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.20	28.35 ng	2433520	1,4-Dichlorobenzene-d4	6.76

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	91
2	Propanenitrile, 3-[(phenylmethyl...	160	C10H12N2	000706-03-6	72
3	Phenethylamine, N-benzyl-p-chloro-	245	C15H16ClN	013622-43-0	72
4	1,2-Ethanediamine, N,N'-bis(phen...	240	C16H20N2	000140-28-3	64
5	Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	56





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122116\
Data File : BF091867.D
Acq On : 22 Dec 2016 6:39
Operator : UM/SJ
Sample : H6156-13
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.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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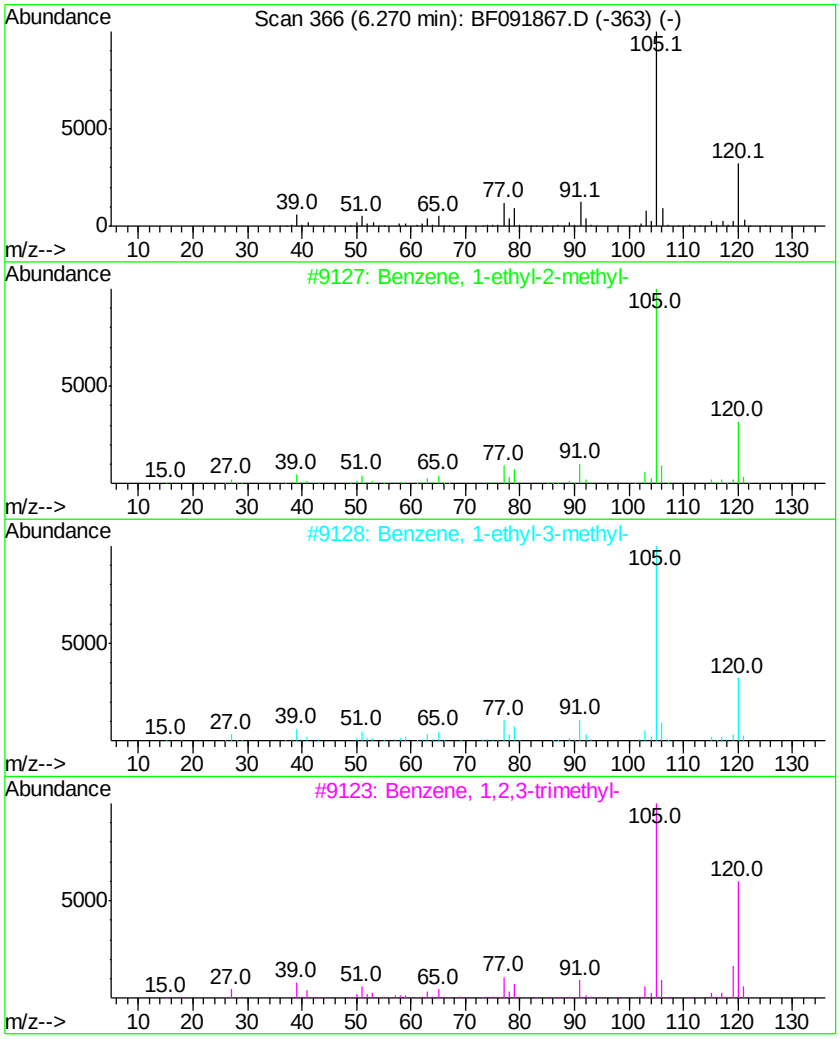
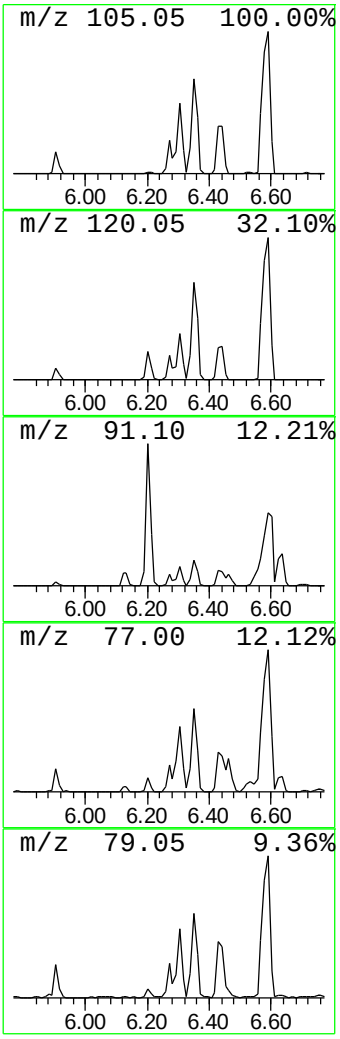
Peak Number 6 Benzene, 1-ethyl-2-methyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.27	23.72 ng	2035850	1,4-Dichlorobenzene-d4	6.76

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, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





Data Path : Z:\HPCHEM1\BNA F\DATA\BF122116\
Data File : BF091867.D
Acq On : 22 Dec 2016 6:39
Operator : UM/SJ
Sample : H6156-13
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.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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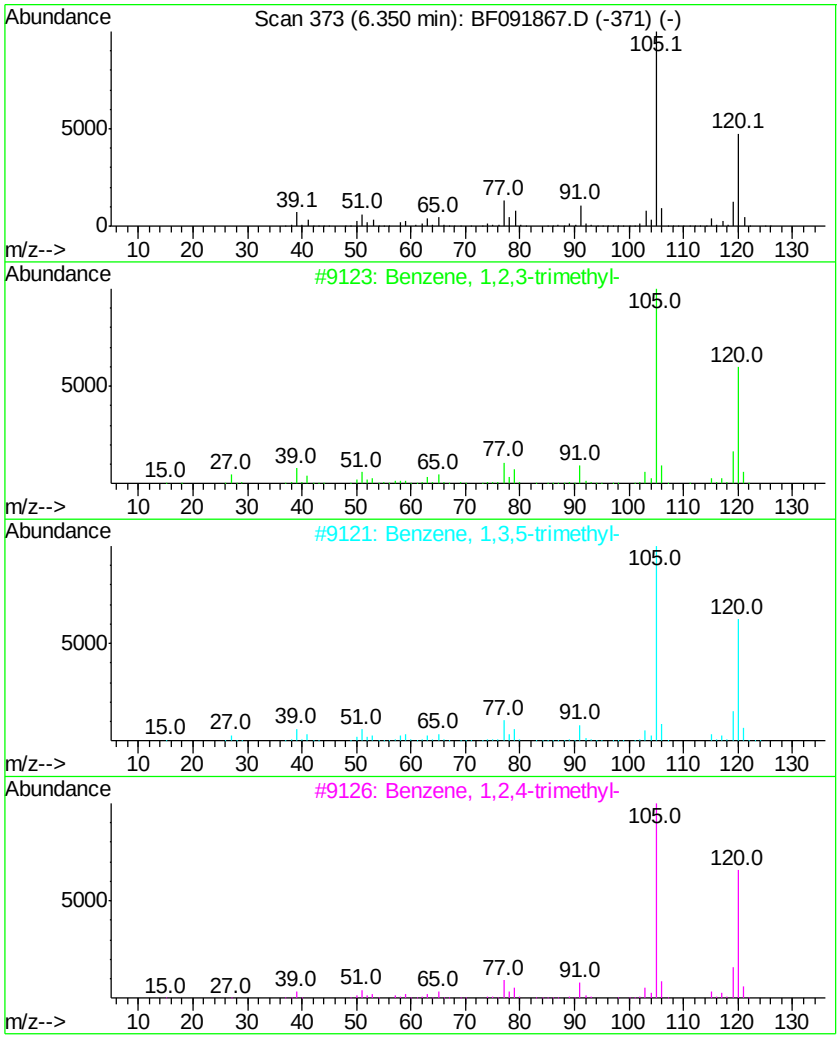
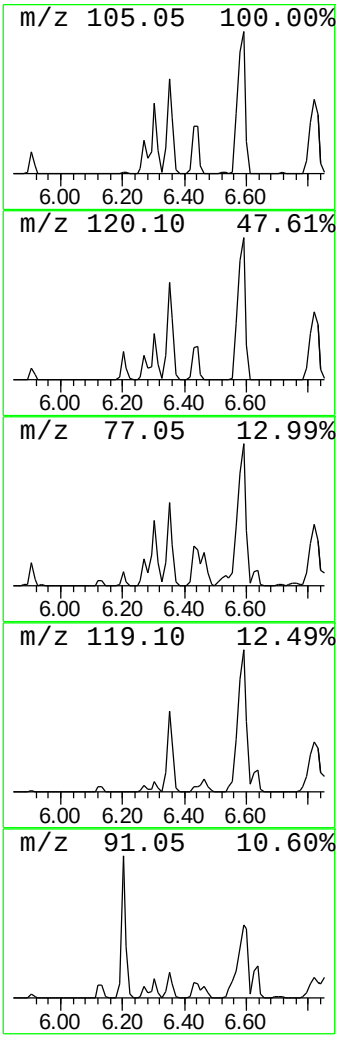
Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.35	87.69 ng	7525610	1,4-Dichlorobenzene-d4	6.76

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2	Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	95
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
4	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90





Data Path : Z:\HPCHEM1\BNA F\DATA\BF122116\
Data File : BF091867.D
Acq On : 22 Dec 2016 6:39
Operator : UM/SJ
Sample : H6156-13
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

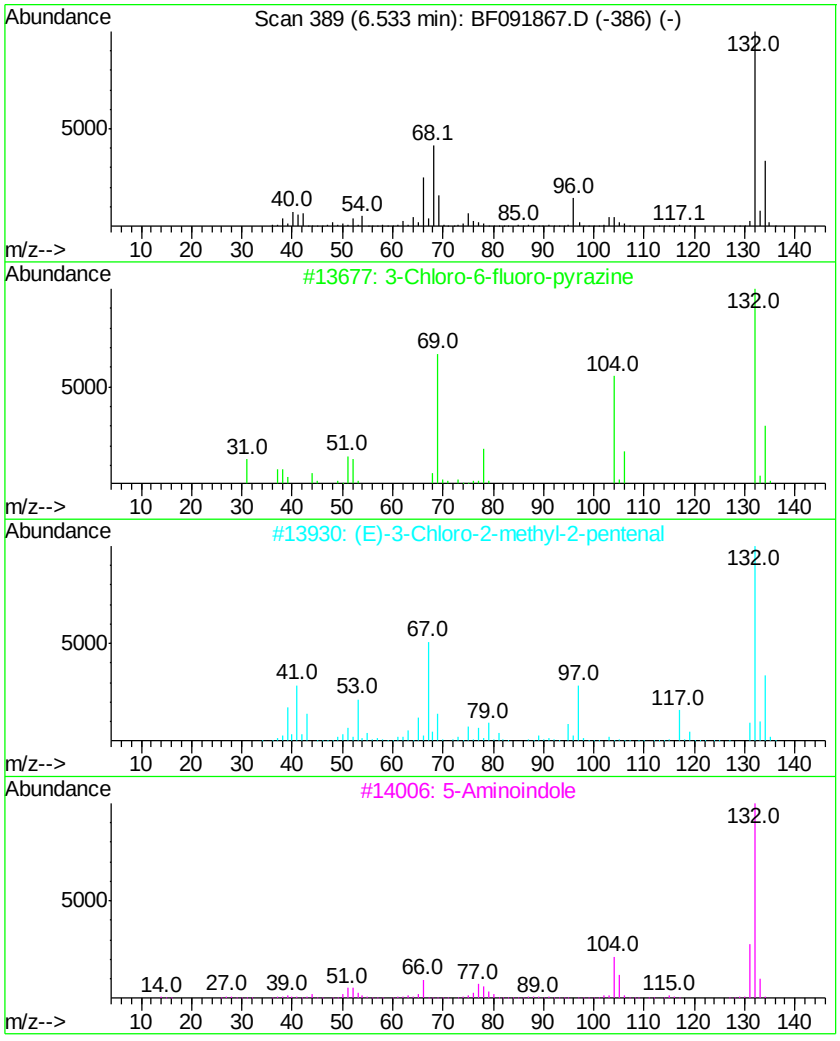
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

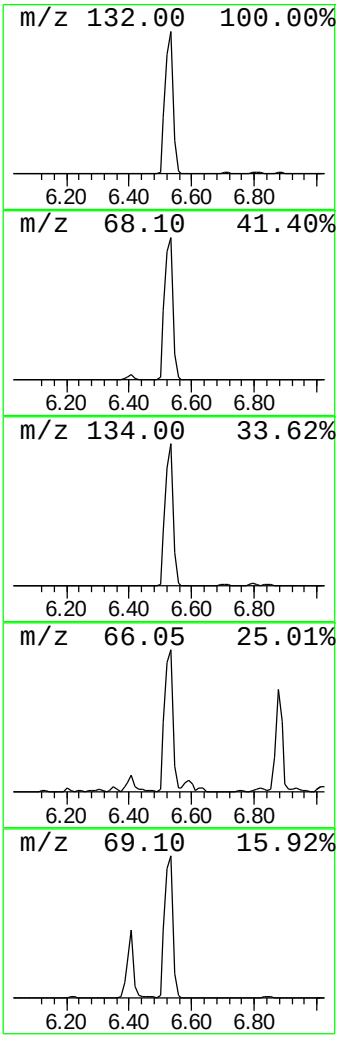
****

Peak Number 9 unknown6.53 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.53	99.68 ng	8555280	1,4-Dichlorobenzene-d4	6.76

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	27
3	5-Aminoindole	132	C8H8N2	005192-03-0	12
4	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5	2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF122116\
Data File : BF091867.D
Acq On : 22 Dec 2016 6:39
Operator : UM/SJ
Sample : H6156-13
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.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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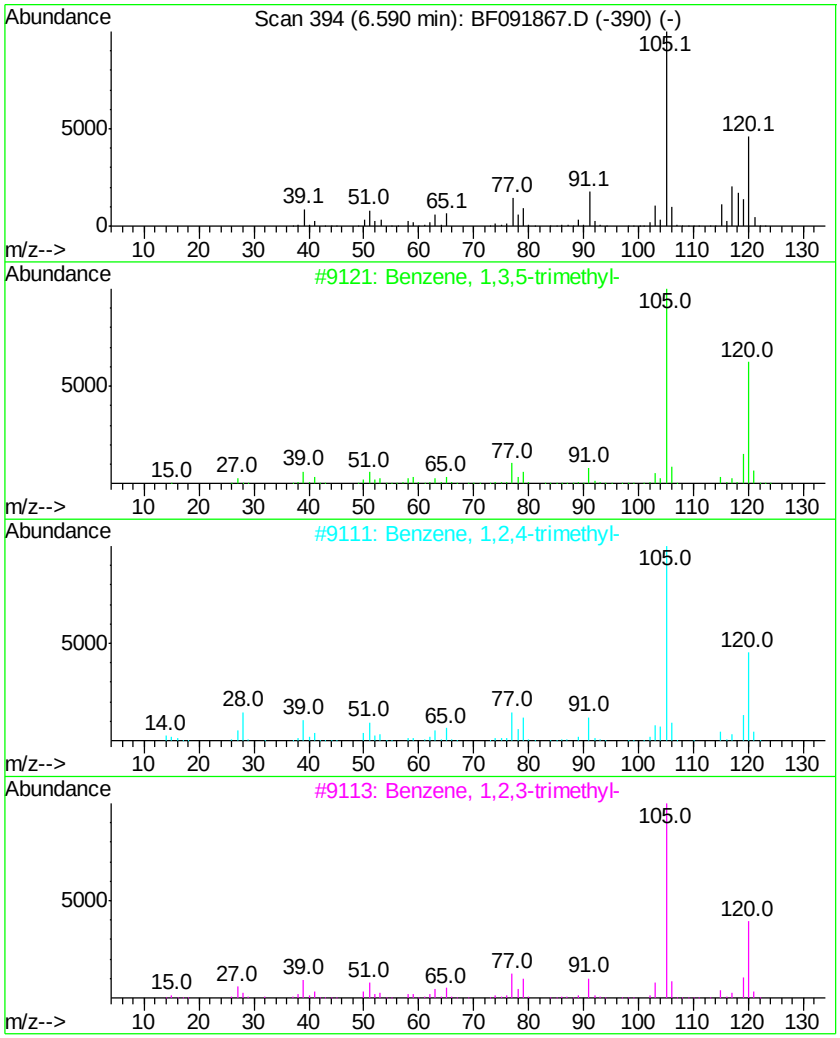
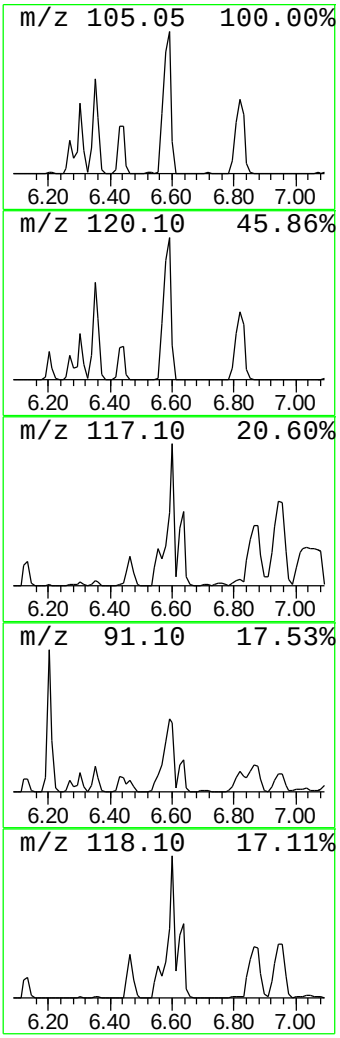
Peak Number 10 Benzene, 1,3,5-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.59	294.27 ng	25255300	1,4-Dichlorobenzene-d4	6.76

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	76
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	76
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	76
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	70
5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	64





Data Path : Z:\HPCHEM1\BNA F\DATA\BF122116\
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Operator : UM/SJ
Sample : H6156-13
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.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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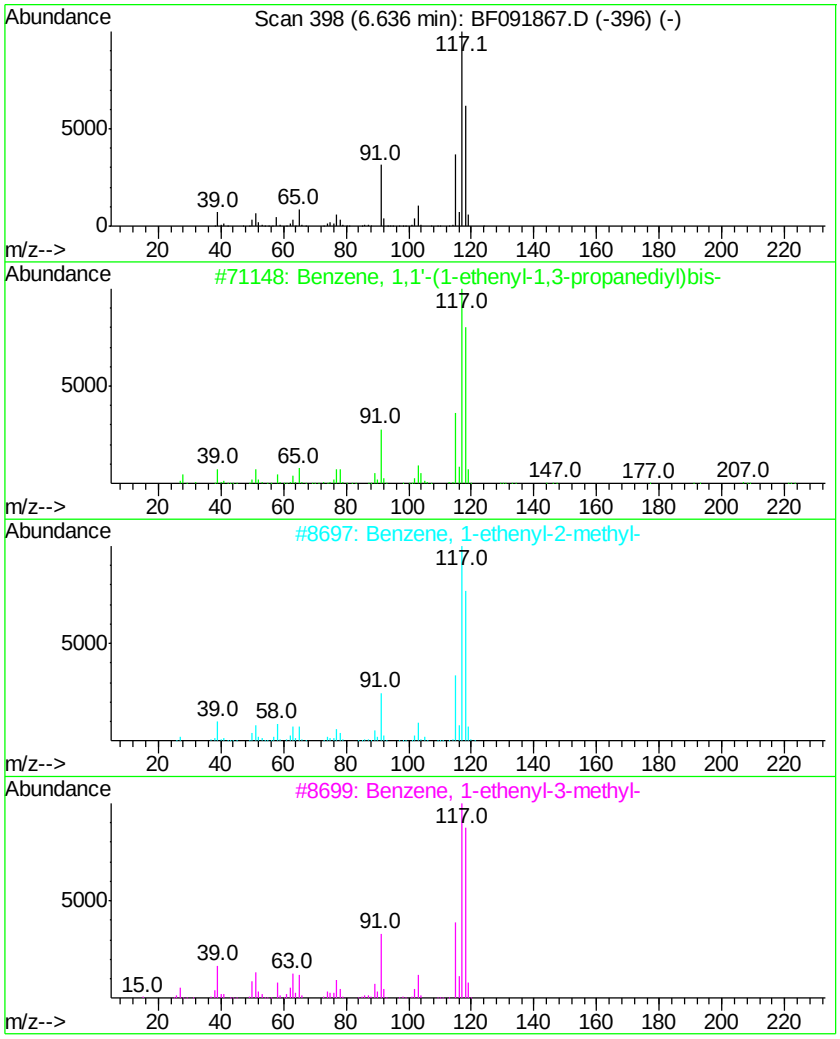
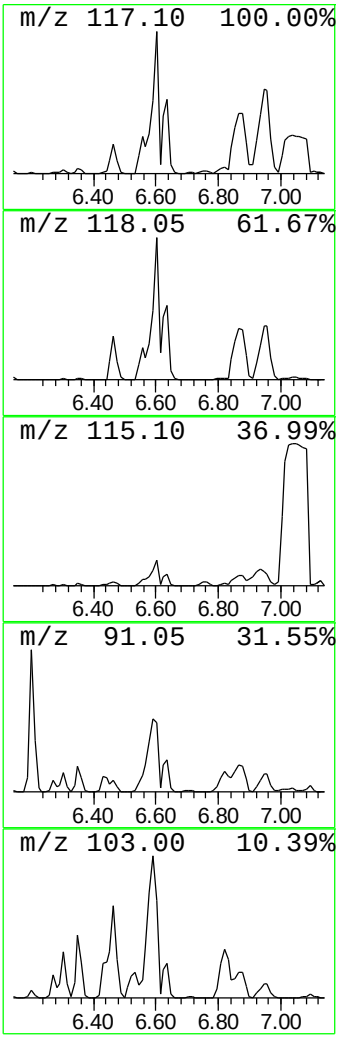
Peak Number 11 Benzene, 1,1'-(1-ethenyl-1,... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.64	50.81 ng	4360950	1,4-Dichlorobenzene-d4	6.76

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,1'-(1-ethenyl-1,3-pro...	222	C17H18	061141-97-7	90
2	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	87
3	Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	87
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	87
5	Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	83





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122116\
Data File : BF091867.D
Acq On : 22 Dec 2016 6:39
Operator : UM/SJ
Sample : H6156-13
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.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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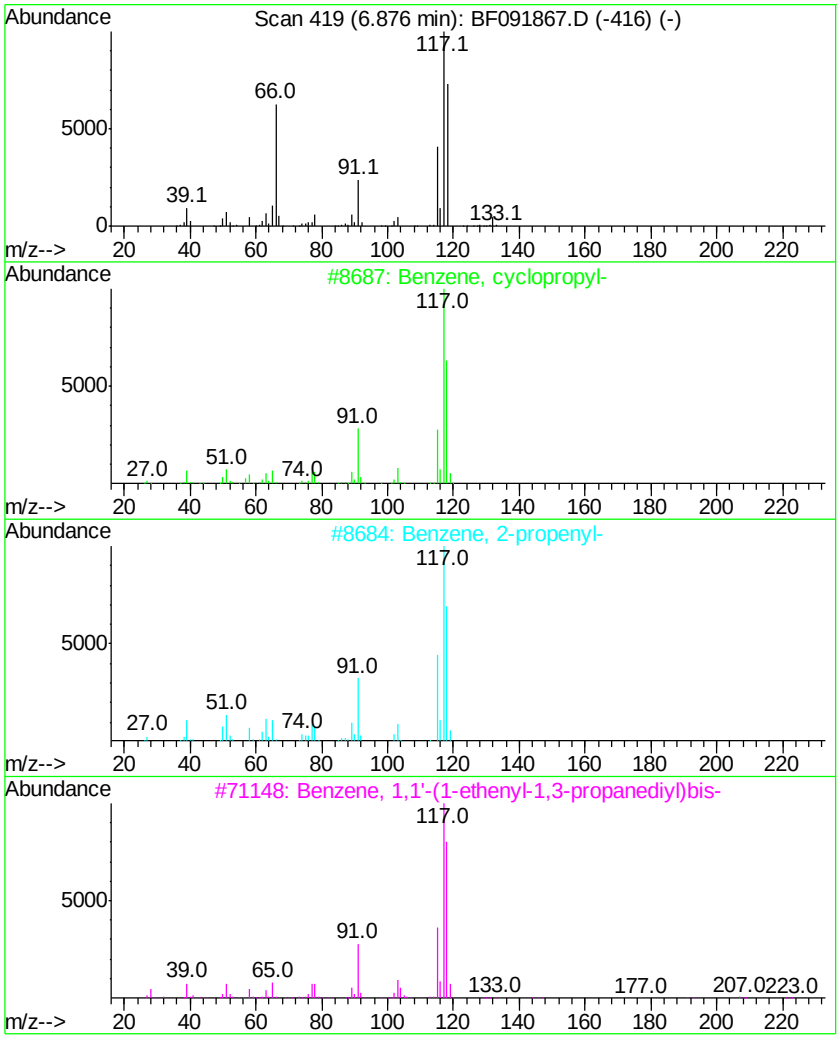
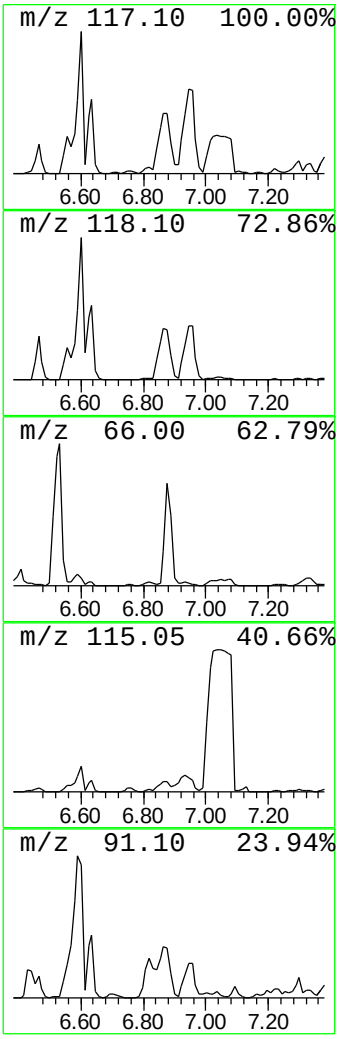
Peak Number 13 unknown6.88 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.88	46.44 ng	3985650	1,4-Dichlorobenzene-d4	6.76

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, cyclopropyl-	118	C9H10	000873-49-4	42
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	38
3	Benzene, 1,1'-(1-ethenyl-1,3-pro...	222	C17H18	061141-97-7	38
4	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	38
5	Benzene, 1-propenyl-	118	C9H10	000637-50-3	38





Data Path : Z:\HPCHEM1\BNA F\DATA\BF122116\
Data File : BF091867.D
Acq On : 22 Dec 2016 6:39
Operator : UM/SJ
Sample : H6156-13
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.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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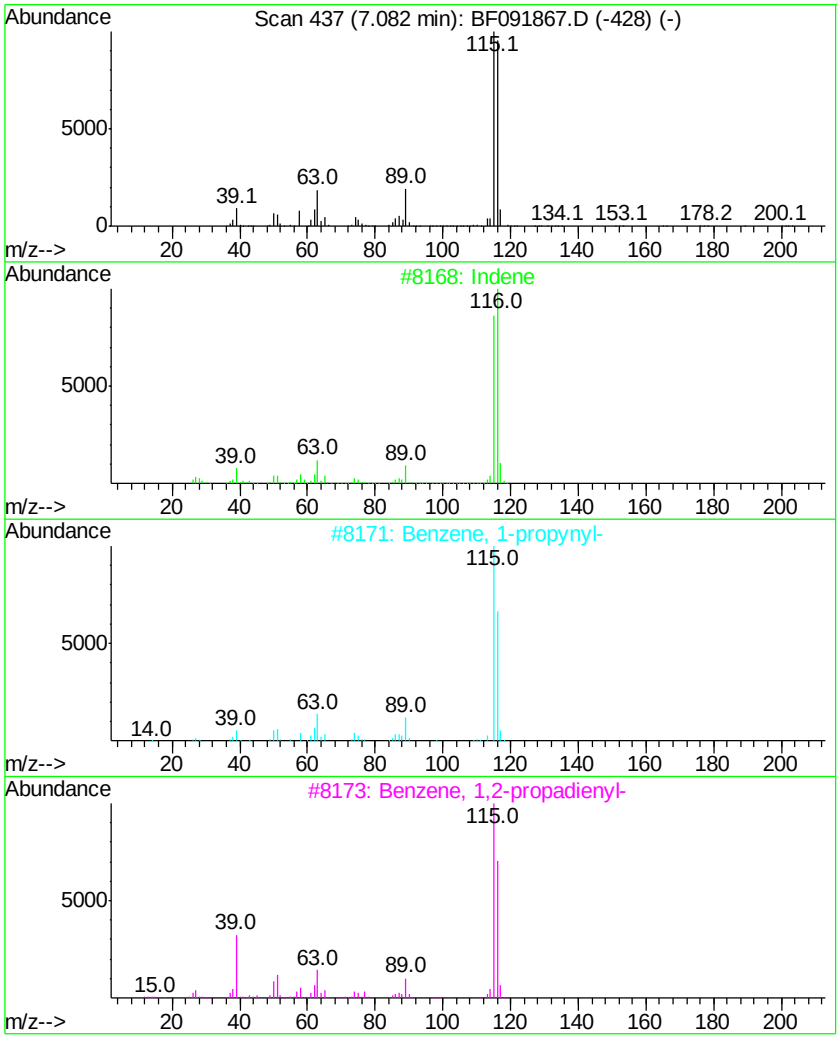
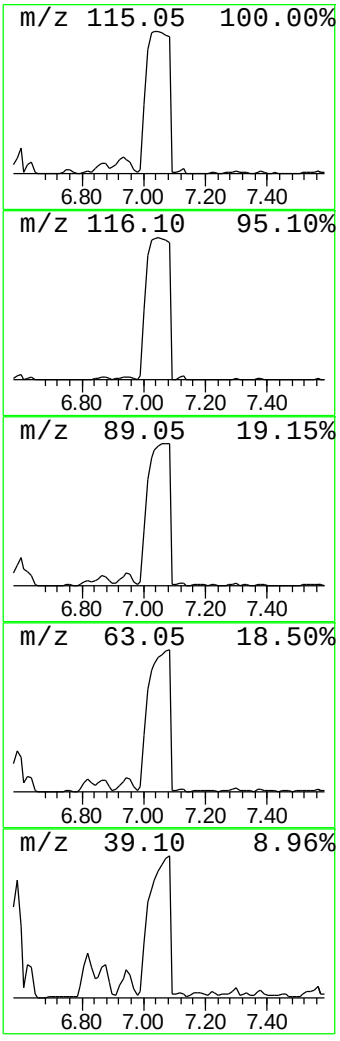
Peak Number 15 Indene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.08	776.65 ng	66655700	1,4-Dichlorobenzene-d4	6.76

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene	116	C9H8	000095-13-6	96
2	Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
3	Benzene, 1,2-propadienyl-	116	C9H8	002327-99-3	91
4	Benzene,1-ethynyl-2-methyl-	116	C9H8	000766-47-2	90
5	Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	90





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122116\
Data File : BF091867.D
Acq On : 22 Dec 2016 6:39
Operator : UM/SJ
Sample : H6156-13
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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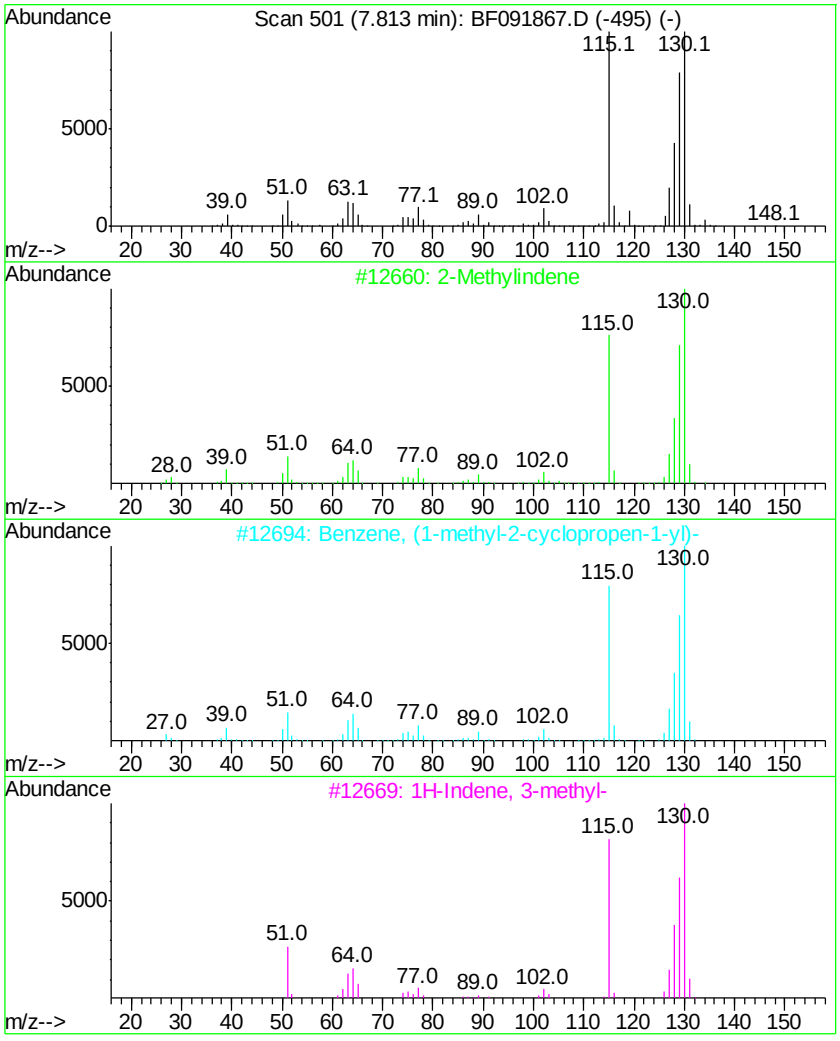
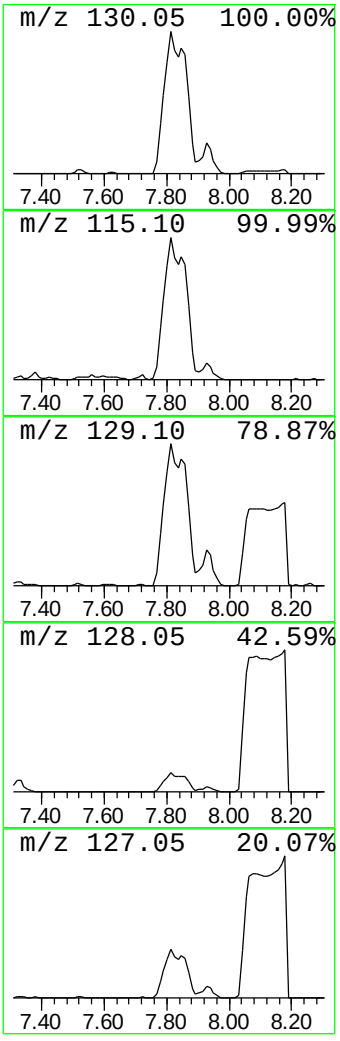
Peak Number 16 2-Methylindene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.81	123.38 ng	28648600	Naphthalene-d8	8.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Methylindene	130	C10H10	002177-47-1	97
2	Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	97
3	1H-Indene, 3-methyl-	130	C10H10	000767-60-2	95
4	1H-Indene, 1-methyl-	130	C10H10	000767-59-9	95
5	1,4-Dihydronaphthalene	130	C10H10	000612-17-9	95





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Data File : BF091867.D
Acq On : 22 Dec 2016 6:39
Operator : UM/SJ
Sample : H6156-13
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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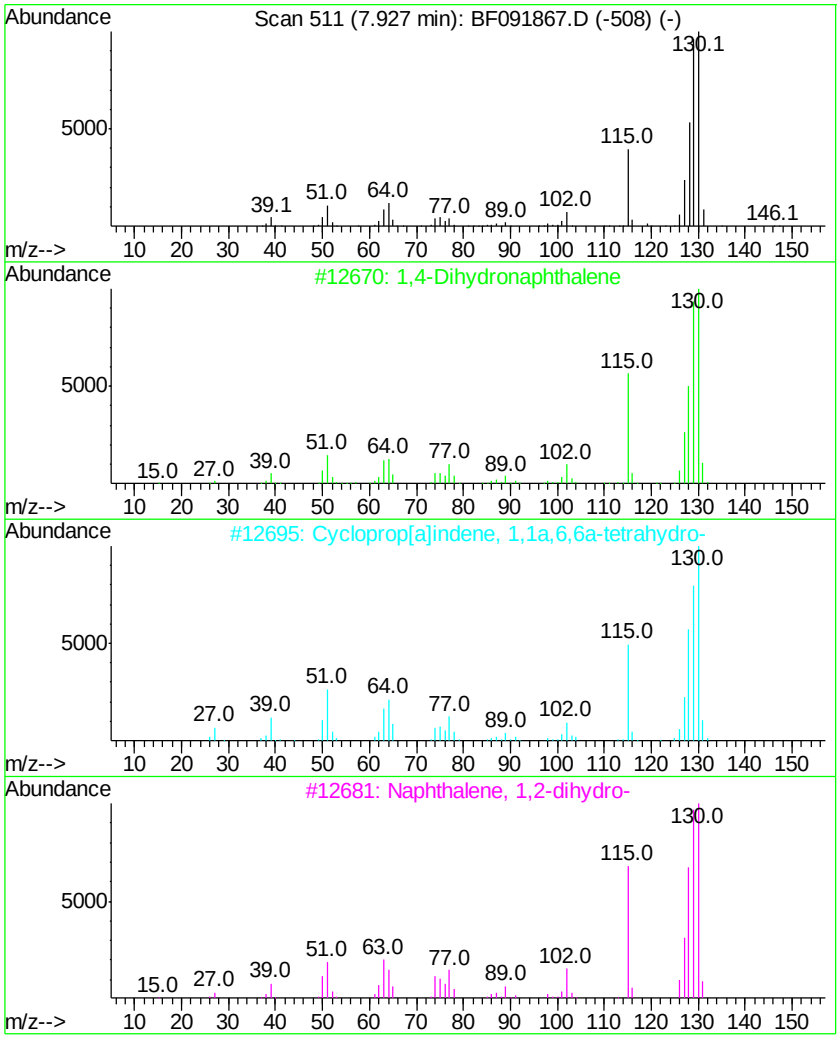
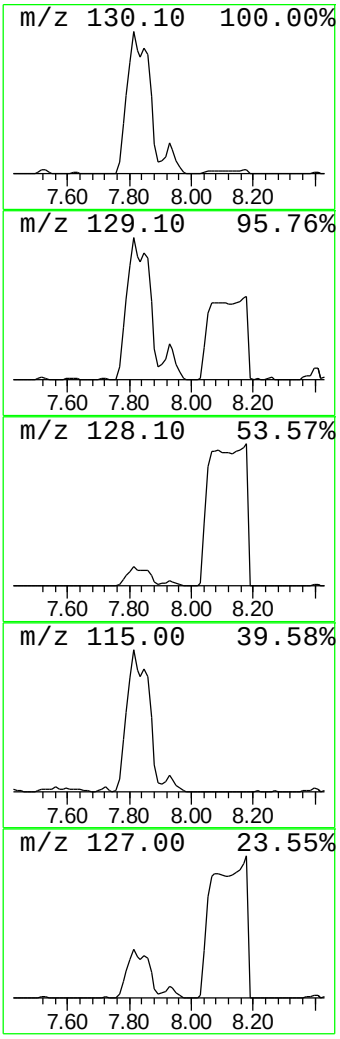
Peak Number 17 1,4-Dihydronaphthalene Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.93	20.00 ng	4643980	Naphthalene-d8	8.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,4-Dihydronaphthalene	130	C10H10	000612-17-9	94
2	Cycloprop[a]indene, 1,1a,6,6a-te...	130	C10H10	015677-15-3	83
3	Naphthalene, 1,2-dihydro-	130	C10H10	000447-53-0	76
4	Benzene, (1-methylene-2-propenyl)-	130	C10H10	002288-18-8	64
5	Triquinacene	130	C10H10	006053-74-3	64





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Operator : UM/SJ
Sample : H6156-13
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.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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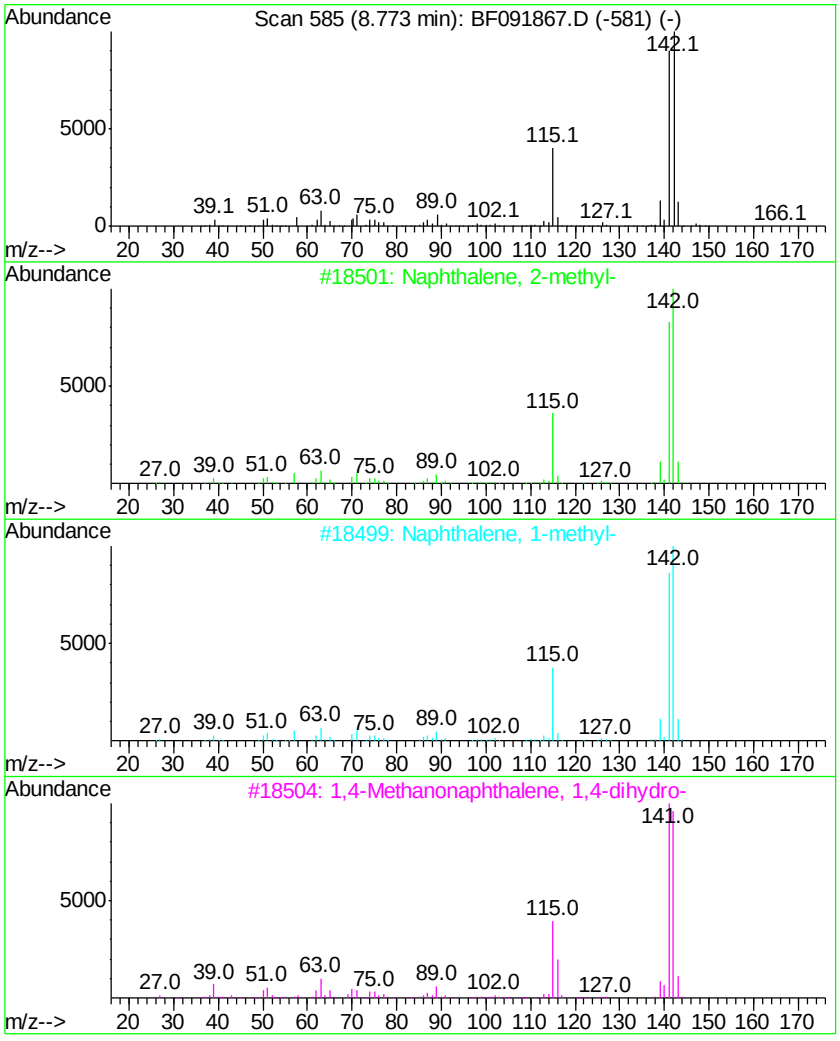
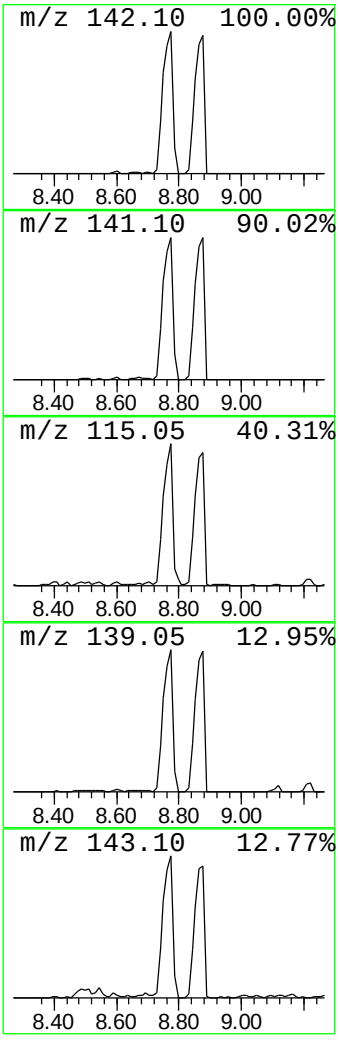
Peak Number 18 Naphthalene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.77	108.13 ng	25108400	Naphthalene-d8	8.03

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94
4	Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	91
5	Benzocycloheptatriene	142	C11H10	000264-09-5	91





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Operator : UM/SJ
Sample : H6156-13
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.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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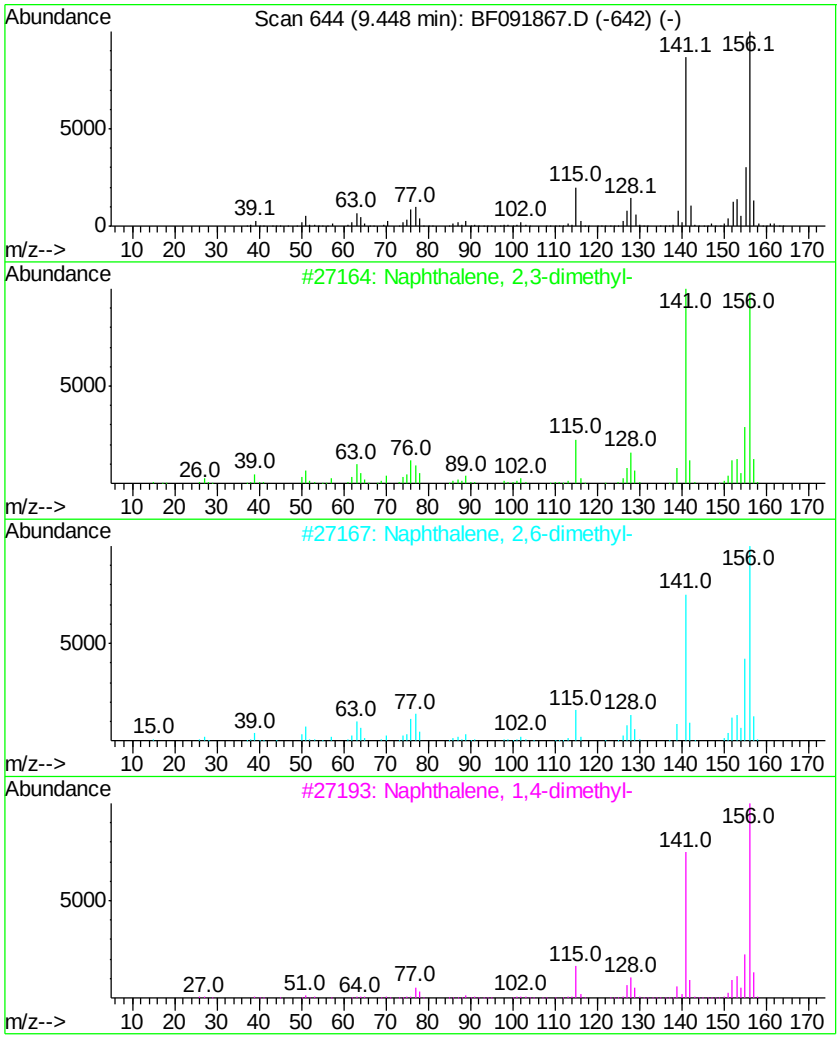
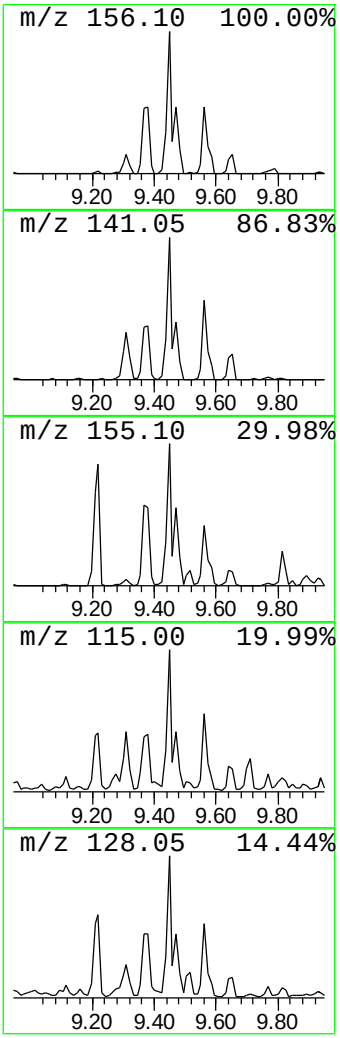
Peak Number 20 Naphthalene, 2,3-dimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.45	19.96 ng	3662060	Acenaphthene-d10	9.79

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
3	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
4	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
5	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96





Tentatively Identified Compound (LSC) summary

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Data File : BF091867.D
Acq On : 22 Dec 2016 6:39
Operator : UM/SJ
Sample : H6156-13
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzene, cyclopro...	6.13	17.5	ng	1501500	1	6.76	1716490	20.0
Benzene, propyl-	6.20	28.4	ng	2433520	1	6.76	1716490	20.0
Benzene, 1-ethyl-...	6.27	23.7	ng	2035850	1	6.76	1716490	20.0
Benzene, 1,2,3-tr...	6.35	87.7	ng	7525610	1	6.76	1716490	20.0
unknown6.53	6.53	99.7	ng	8555280	1	6.76	1716490	20.0
Benzene, 1,3,5-tr...	6.59	294.3	ng	25255300	1	6.76	1716490	20.0
Benzene, 1,1'-(1-...	6.64	50.8	ng	4360950	1	6.76	1716490	20.0
unknown6.88	6.88	46.4	ng	3985650	1	6.76	1716490	20.0
Indene	7.08	776.6	ng	66655700	1	6.76	1716490	20.0
2-Methylindene	7.81	123.4	ng	28648600	2	8.03	4643980	20.0
1,4-Dihydronaphth...	7.93	20.0	ng	4643980	2	8.03	4643980	20.0
Naphthalene, 1-me...	8.77	108.1	ng	25108400	2	8.03	4643980	20.0
Naphthalene, 2,3-...	9.45	20.0	ng	3662060	3	9.79	3668550	20.0