LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060416\
Data File : BF087696.D
Acq On : 5 Jun 2016 8:41
Operator : UM/SJ
Sample : H3345-01
Misc :
ALS Vial : 38 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-BF060416.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	1.770	15	16	20	rVV	314409	406878	2.07%	0.224%
2	1.850	20	23	26	rVV	794020	1026354	5.21%	0.566%
3	1.907	26	28	43	rVB	2396951	4129922	20.97%	2.279%
4	3.896	192	202	203	rBV	5048535	19691576	100.00%	10.865%
5	5.324	324	327	330	rBV	3453346	4102633	20.83%	2.264%
6	5.473	334	340	342	rBV2	4891102	12393755	62.94%	6.838%
7	5.724	356	362	365	rVV3	5338290	17162290	87.16%	9.469%
8	6.022	386	388	390	rVB	332838	278567	1.41%	0.154%
9	6.239	405	407	409	rBV	335445	344550	1.75%	0.190%
10	6.319	412	414	416	rBV	770630	656940	3.34%	0.362%
11	6.387	418	420	421	rVV	506779	489829	2.49%	0.270%
12	6.422	421	423	425	rVV	1173494	1313825	6.67%	0.725%
13	6.467	425	427	431	rVV2	2393326	3285515	16.68%	1.813%
14	6.547	431	434	435	rVV	1056860	1106720	5.62%	0.611%
15	6.582	435	437	438	rVV	746270	1029300	5.23%	0.568%
16	6.627	438	441	443	rVV	2269797	2950149	14.98%	1.628%
17	6.696	443	447	449	rVV2	4403638	8650539	43.93%	4.773%
18	6.742	449	451	453	rVV	1758338	1693995	8.60%	0.935%
19	6.867	460	462	464	rVV	650575	766100	3.89%	0.423%
20	6.924	464	467	469	rVV	1764358	2653879	13.48%	1.464%
21	6.970	469	471	473	rVV	1140677	2007386	10.19%	1.108%
22	7.027	473	476	481	rVB2	1814724	4926457	25.02%	2.718%
23	7.222	491	493	495	rVB2	146062	266020	1.35%	0.147%
24	7.267	495	497	500	rBV	379809	439461	2.23%	0.242%
25	7.336	500	503	504	rVV2	251151	294113	1.49%	0.162%
26	7.427	506	511	514	rVV2	1624466	3207098	16.29%	1.770%
27	7.485	514	516	518	rVV	510872	634134	3.22%	0.350%
28	7.565	521	523	525	rVV2	127053	217022	1.10%	0.120%
29	7.645	525	530	531	rVV3	235130	605192	3.07%	0.334%
30	7.667	531	532	534	rVV	663239	582507	2.96%	0.321%
31	7.702	534	535	537	rVV	336021	350963	1.78%	0.194%
32	7.736	537	538	543	rVV	114301	222254	1.13%	0.123%
33	7.827	543	546	549	rVV	276677	339893	1.73%	0.188%
34	7.907	549	553	555	rVV2	3507613	7910191	40.17%	4.364%

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Acq On : 5 Jun 2016 8:41
Operator : UM/SJ
Sample : H3345-01
Misc :
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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-BF060416.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	7.953	555	557	560	rVV	3289310	5948464	30.21%	3.282%
36	8.033	560	564	567	rVB2	863416	1467547	7.45%	0.810%
37	8.182	572	577	578	rBV	4957278	13043780	66.24%	7.197%
38	8.307	586	588	591	rBV	1125744	1133529	5.76%	0.625%
39	8.513	600	606	607	rBV2	164880	389229	1.98%	0.215%
40	8.593	610	613	614	rVV2	340503	679392	3.45%	0.375%
41	8.650	616	618	620	rVV	314430	569465	2.89%	0.314%
42	8.696	620	622	624	rVV2	274281	507288	2.58%	0.280%
43	8.742	624	626	628	rVV	498457	704982	3.58%	0.389%
44	8.822	628	633	634	rVV2	374369	824351	4.19%	0.455%
45	8.879	634	638	640	rVV	4831852	9377764	47.62%	5.174%
46	8.982	643	647	649	rVV	4771266	9140241	46.42%	5.043%
47	9.233	665	669	670	rBV	3134348	4277164	21.72%	2.360%
48	9.325	674	677	679	rVB	2235287	1973981	10.02%	1.089%
49	9.416	684	685	688	rVB	332554	412213	2.09%	0.227%
50	9.485	688	691	693	rBV	710363	990057	5.03%	0.546%
51	9.565	695	698	699	rBV	1129512	1495349	7.59%	0.825%
52	9.588	699	700	702	rVV2	709326	619256	3.14%	0.342%
53	9.622	702	703	705	rVB2	560141	453748	2.30%	0.250%
54	9.679	705	708	712	rVB	679190	827476	4.20%	0.457%
55	9.770	712	716	718	rVB	3106600	4200746	21.33%	2.318%
56	9.896	723	727	729	rBV3	1018105	1406380	7.14%	0.776%
57	9.930	729	730	732	rVB	877529	672167	3.41%	0.371%
58	10.102	742	745	747	rVV	245513	241860	1.23%	0.133%
59	10.205	751	754	756	rVV4	115859	243294	1.24%	0.134%
60	10.376	768	769	771	rVV	183984	255328	1.30%	0.141%
61	10.411	771	772	773	rVV	253404	239199	1.21%	0.132%
62	10.445	773	775	778	rVB	419537	434166	2.20%	0.240%
63	10.685	792	796	799	rVV2	2131025	2602562	13.22%	1.436%
64	10.811	802	807	811	rBV	359956	405943	2.06%	0.224%
65	11.108	831	833	837	rBV	316539	398543	2.02%	0.220%
66	11.279	842	848	850	rBV2	133489	282864	1.44%	0.156%
67	11.382	854	857	858	rVV	1253807	1394589	7.08%	0.769%
68	11.405	858	859	861	rVV	2061695	1561620	7.93%	0.862%
69	11.451	861	863	865	rVV	215236	239521	1.22%	0.132%
70	11.611	874	877	878	rBV	173160	200175	1.02%	0.110%
71	11.908	902	903	906	rVV2	214024	239072	1.21%	0.132%

LSC Area Percent Report

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Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_F
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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060416.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

72	12.799	978	981	986	rBV2	114467	210243	1.07%	0.116%
73	12.971	993	996	998	rBV	3214482	3337246	16.95%	1.841%
74	14.011	1084	1087	1089	rBV	1110790	990571	5.03%	0.547%
75	15.440	1209	1212	1215	rBV	589813	714643	3.63%	0.394%

Sum of corrected areas: 181242015

LSC Report - Integrated Chromatogram

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Operator : UM/SJ
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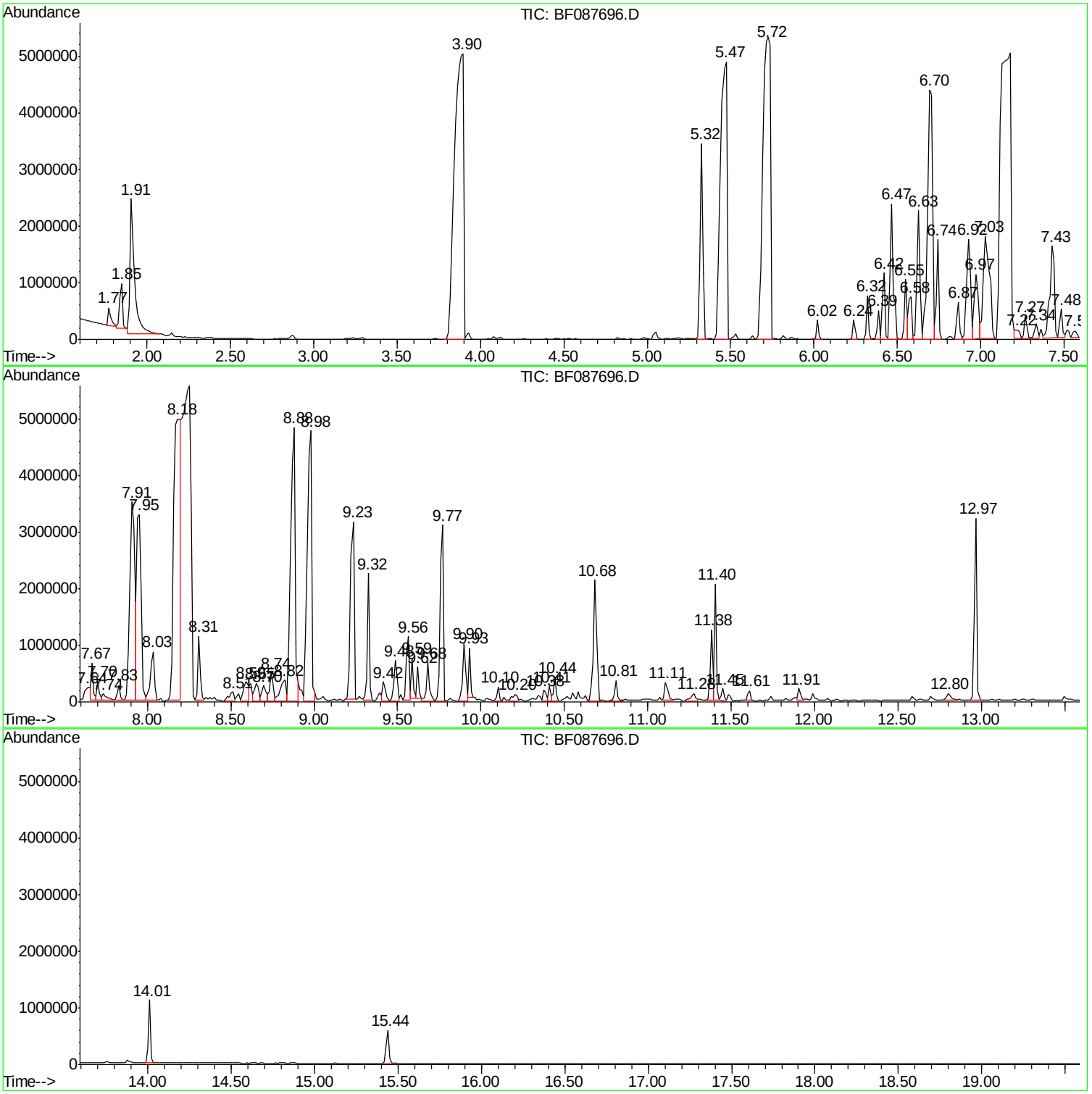
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060416\
Data File : BF087696.D
Acq On : 5 Jun 2016 8:41
Operator : UM/SJ
Sample : H3345-01
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

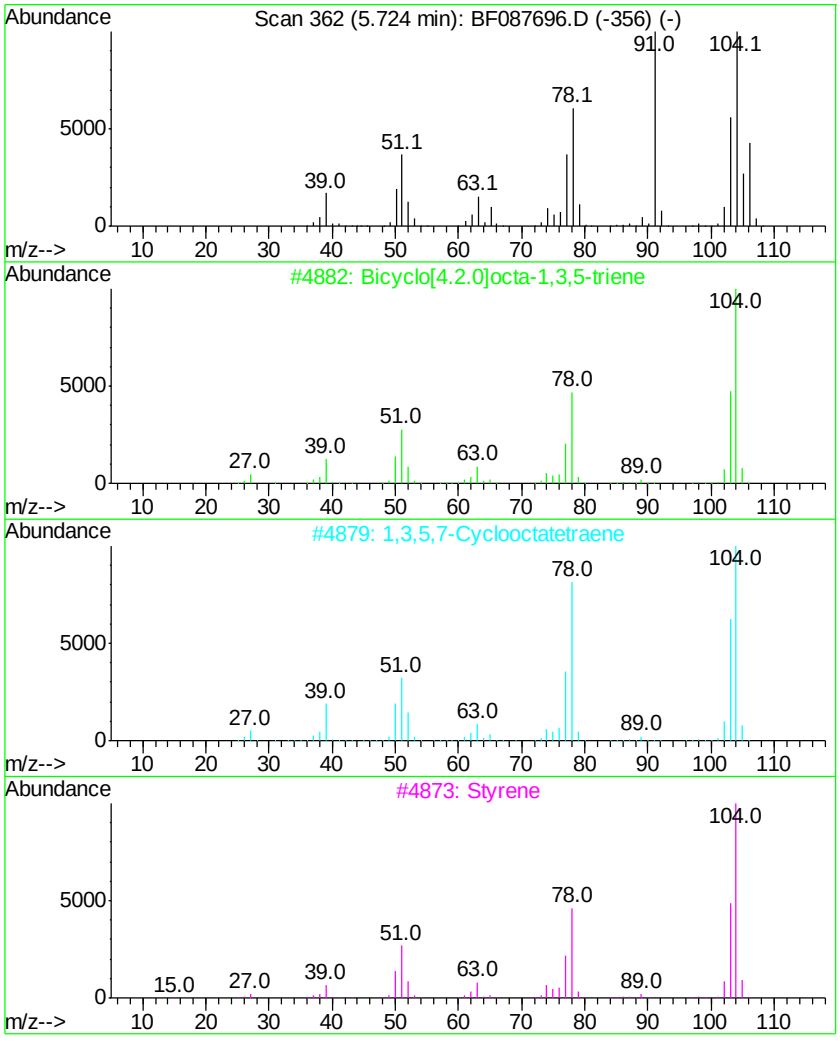
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

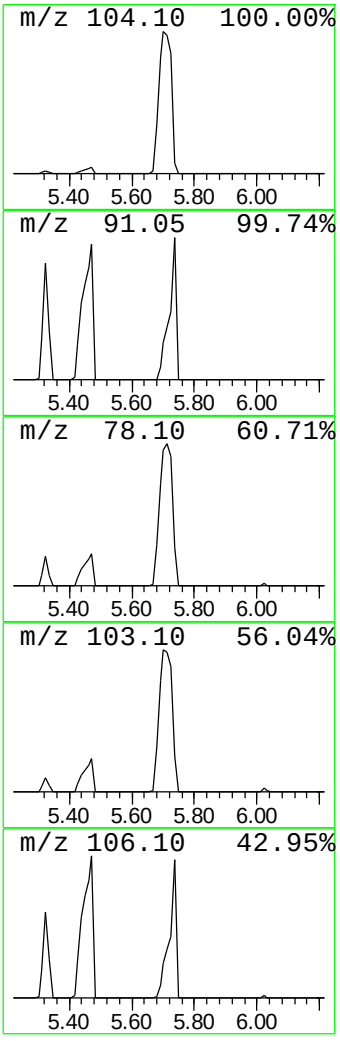
****

Peak Number 5 Bicyclo[4.2.0]octa-1,3,5-tr... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.72	448.04 ng	17162300	1,4-Dichlorobenzene-d4	6.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[4.2.0]octa-1,3,5-triene	104	C8H8	000694-87-1	95
2	1,3,5,7-Cyclooctatetraene	104	C8H8	000629-20-9	94
3	Styrene	104	C8H8	000100-42-5	93
4	1-Cyclohexene, 1-ethynyl-	106	C8H10	1000327-56-5	55
5	1,3,7-Octatrien-5-yne	104	C8H8	016607-77-5	53





Data Path : Z:\HPCHEM1\BNA F\DATA\BF060416\
Data File : BF087696.D
Acq On : 5 Jun 2016 8:41
Operator : UM/SJ
Sample : H3345-01
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

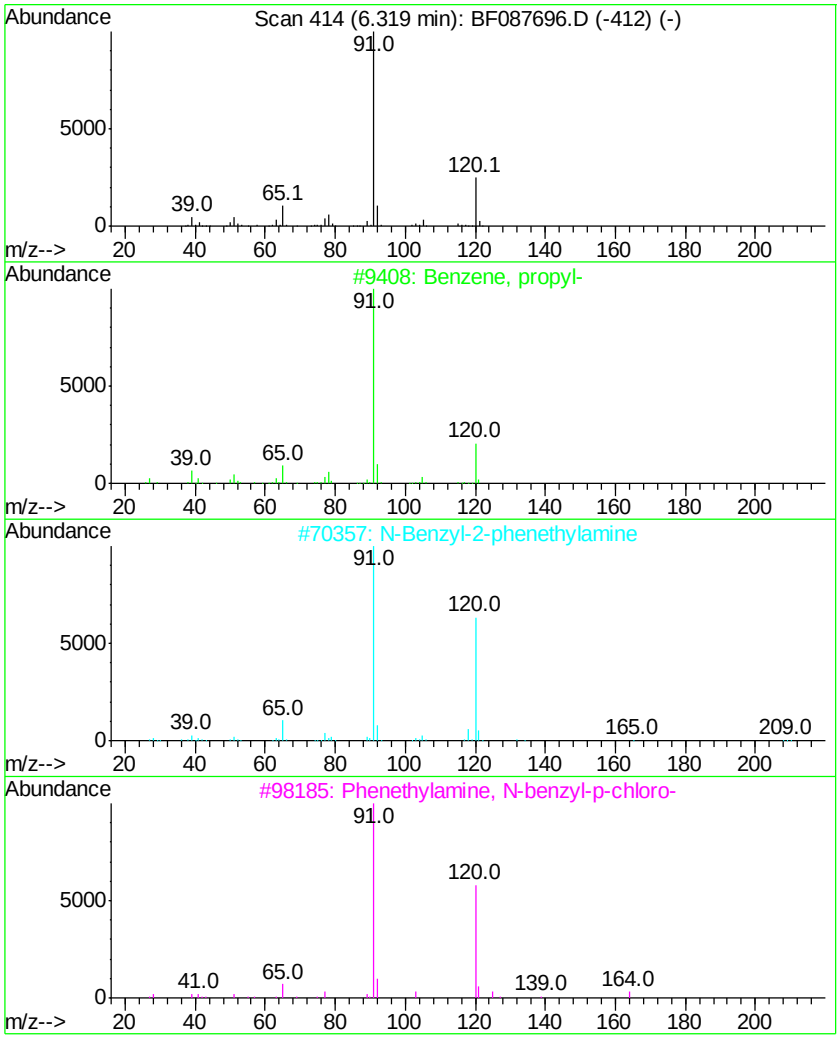
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

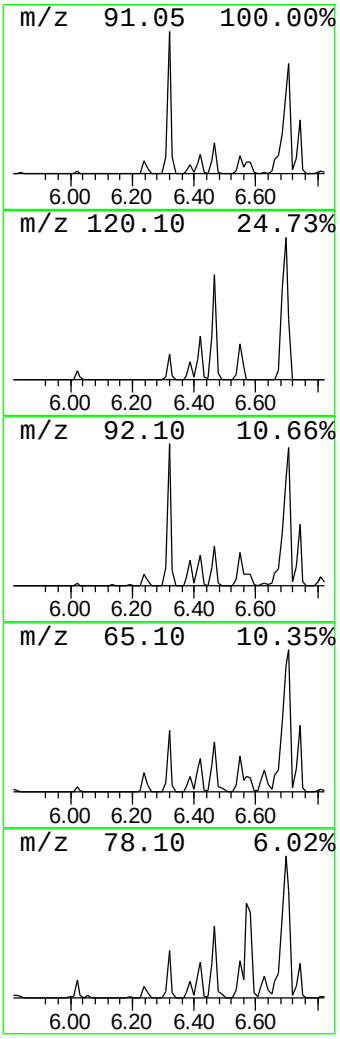
****

Peak Number 6 Benzene, propyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.32	17.15 ng	656940	1,4-Dichlorobenzene-d4	6.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	91
2	N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	74
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	Propanenitrile, 3-[(phenylmethyl...	160	C10H12N2	000706-03-6	59





Data Path : Z:\HPCHEM1\BNA F\DATA\BF060416\
Data File : BF087696.D
Acq On : 5 Jun 2016 8:41
Operator : UM/SJ
Sample : H3345-01
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

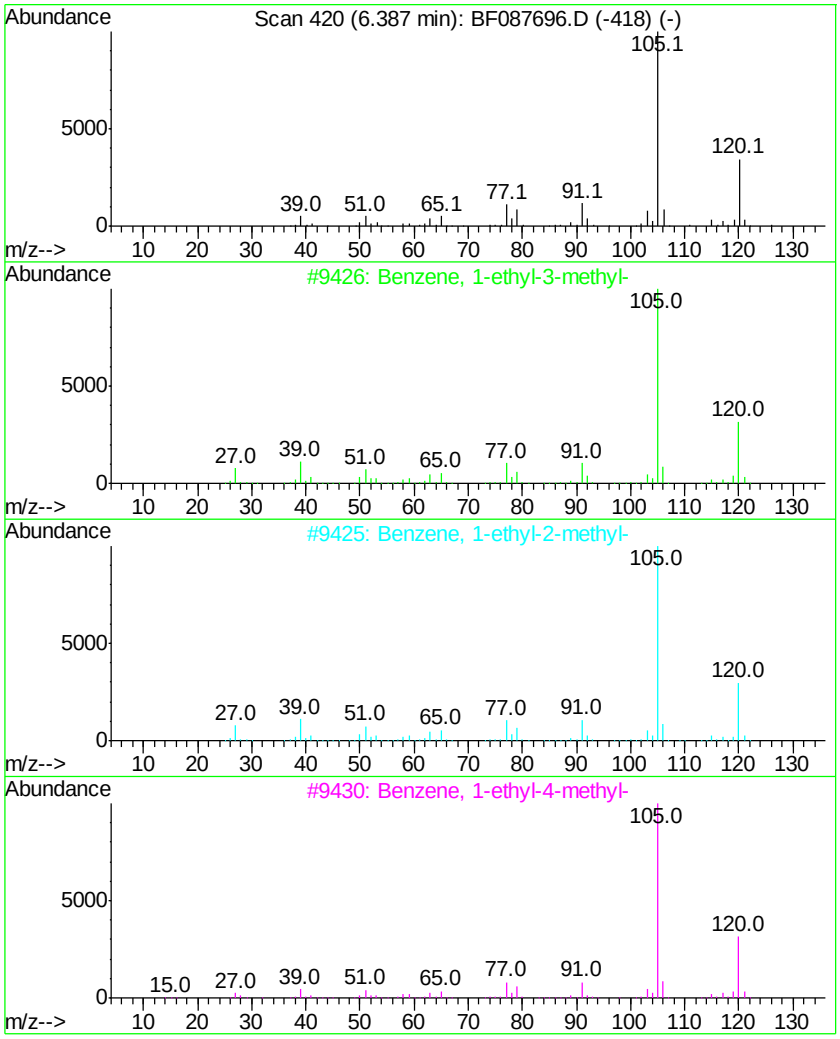
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

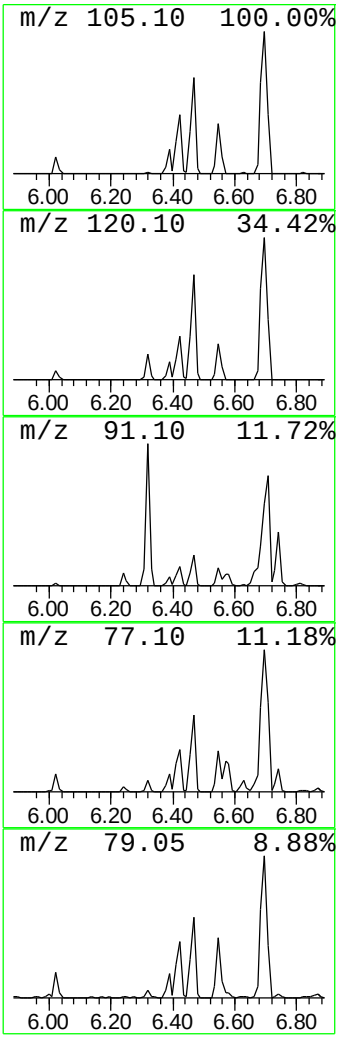
****

Peak Number 7 Benzene, 1-ethyl-3-methyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.39	12.79 ng	489829	1,4-Dichlorobenzene-d4	6.87

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-2-methyl-	120	C9H12	000611-14-3	95
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4	Mesitylene	120	C9H12	000108-67-8	91
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91





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Sample : H3345-01
Misc :
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Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

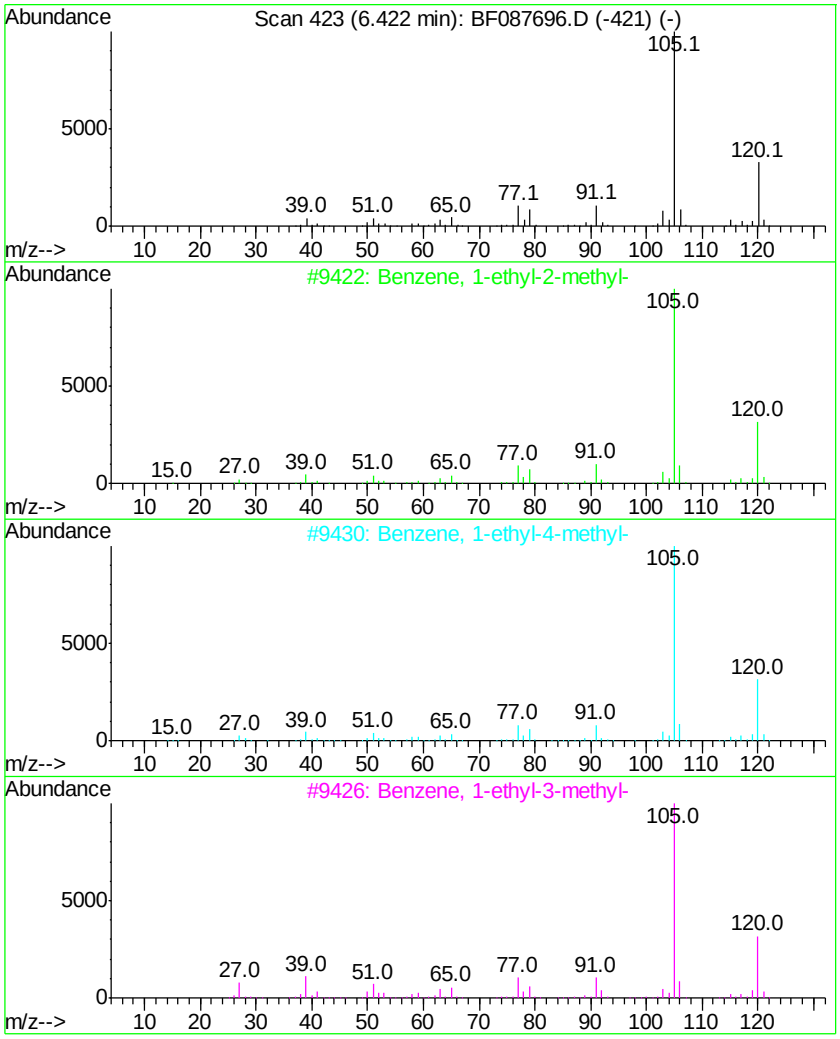
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

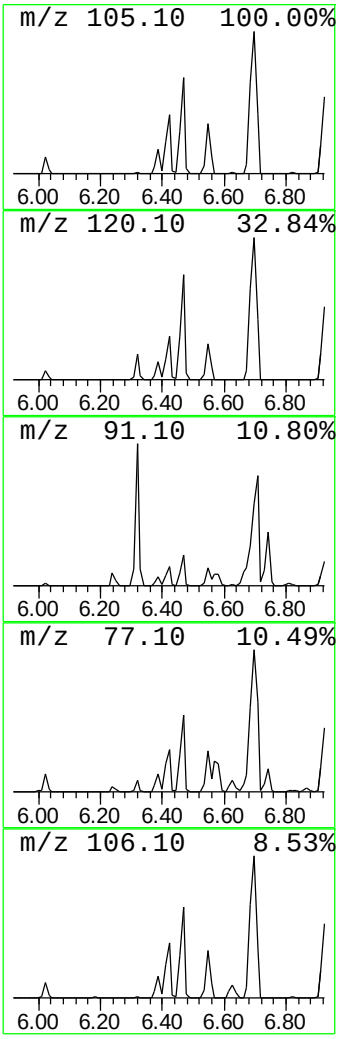
****

Peak Number 8 Benzene, 1-ethyl-2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.42	34.30 ng	1313830	1,4-Dichlorobenzene-d4	6.87

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-4-methyl-	120	C9H12	000622-96-8	91
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91





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Operator : UM/SJ
Sample : H3345-01
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

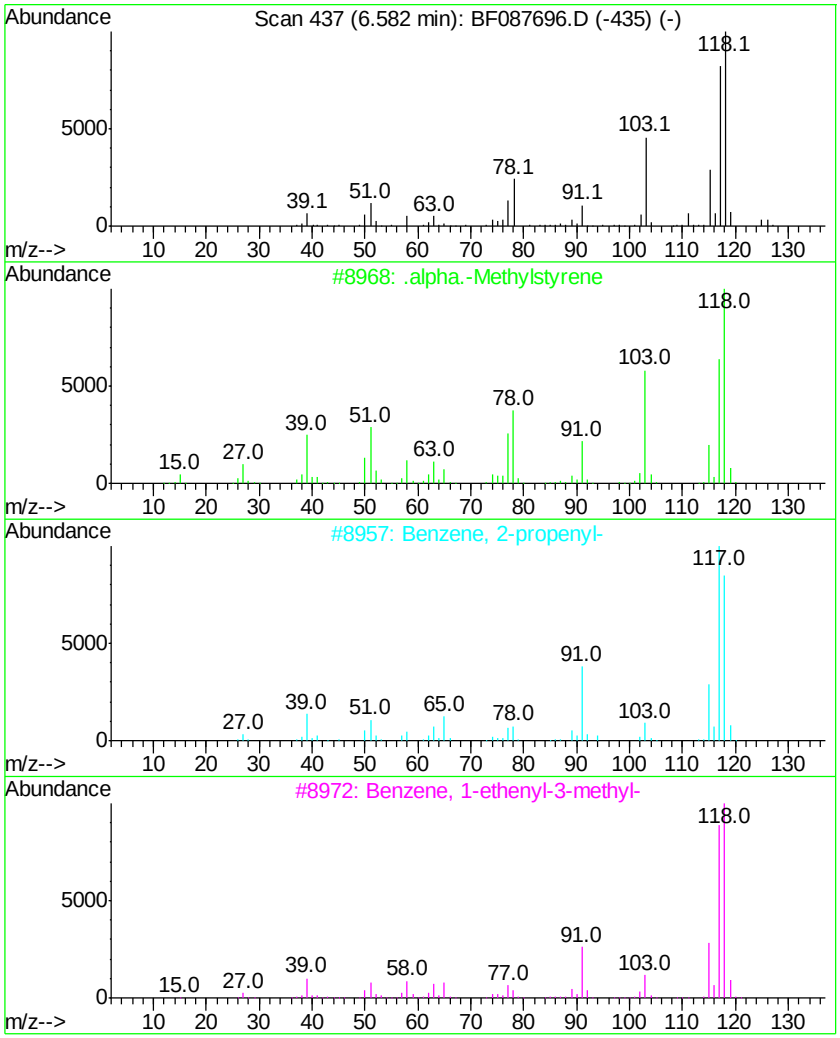
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

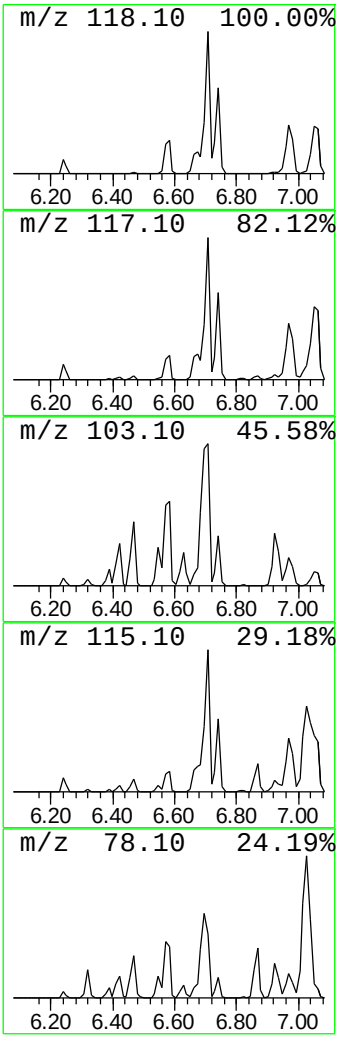
****

Peak Number 10 .alpha.-Methylstyrene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	26.87 ng	1029300	1,4-Dichlorobenzene-d4	6.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.alpha.-Methylstyrene	118	C9H10	000098-83-9	91
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	52
3	Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	40
4	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	40
5	Benzene, 1,1'-(1-ethenyl-1,3-pro...	222	C17H18	061141-97-7	38





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Acq On : 5 Jun 2016 8:41
Operator : UM/SJ
Sample : H3345-01
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

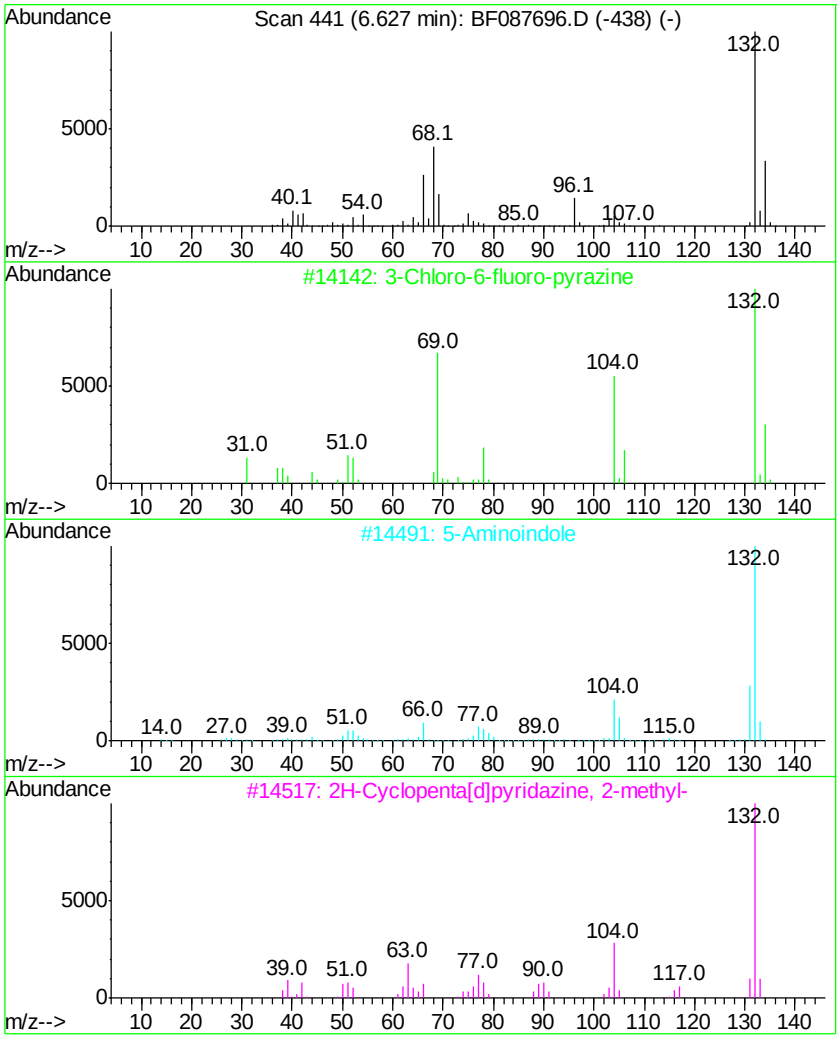
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

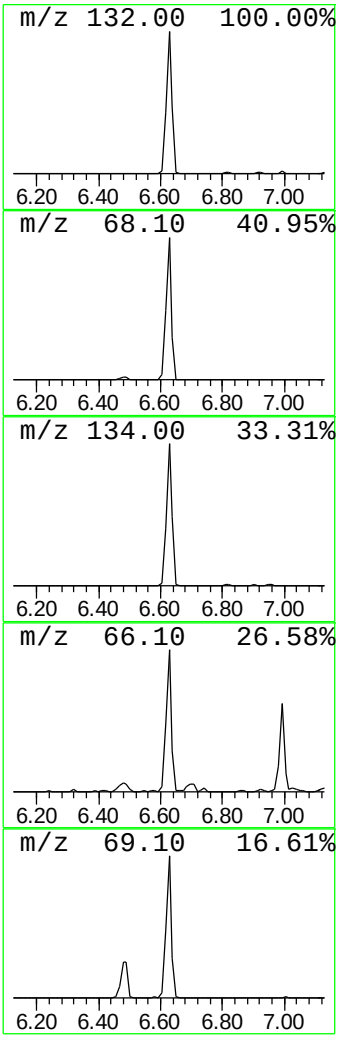
****

Peak Number 11 unknown6.63 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.63	77.02 ng	2950150	1,4-Dichlorobenzene-d4	6.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	5-Aminoindole	132	C8H8N2	005192-03-0	12
3	2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
4	1-Aminoindan	133	C9H11N	034698-41-4	9
5	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF060416\
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Sample : H3345-01
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_F
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

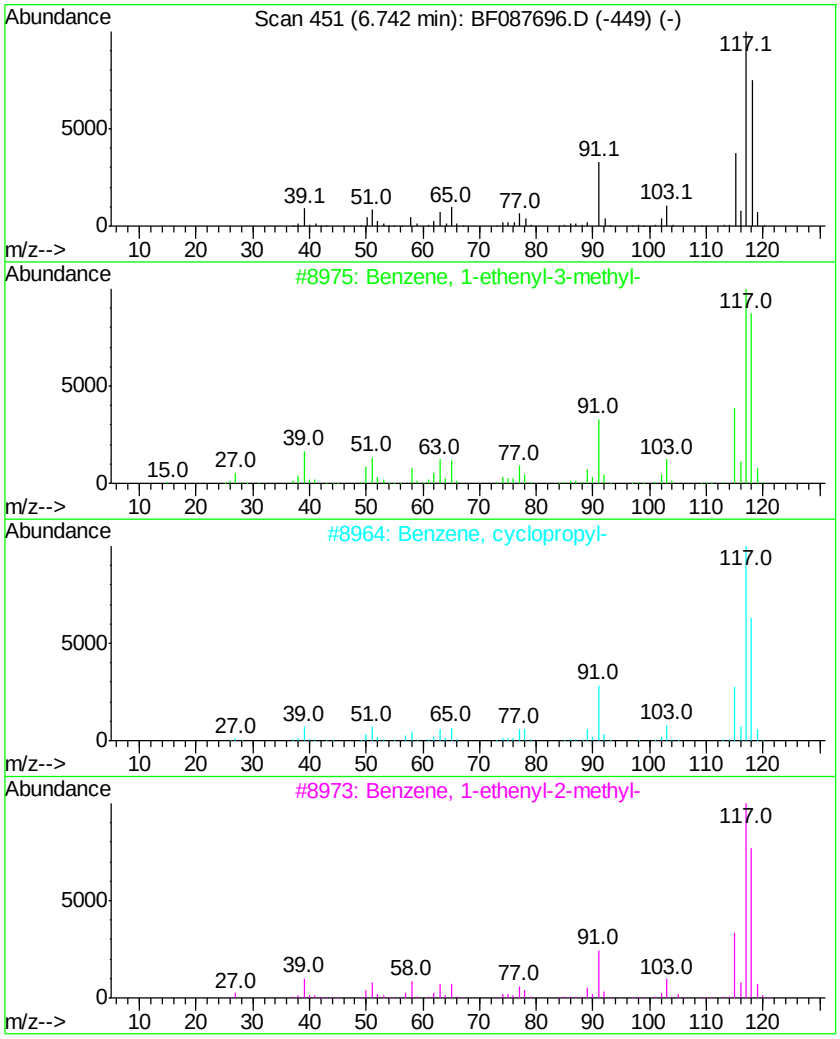
TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

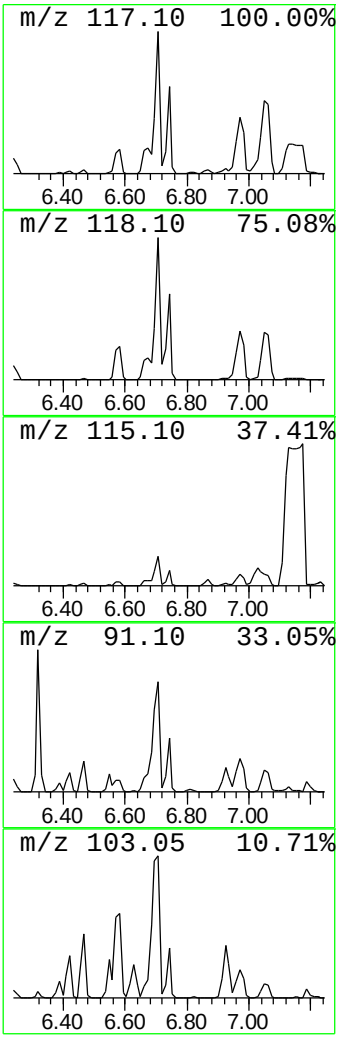
****

Peak Number 13 Benzene, 1-ethenyl-3-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.74	44.22 ng	1694000	1,4-Dichlorobenzene-d4	6.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	93
2	Benzene, cyclopropyl-	118	C9H10	000873-49-4	90
3	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	90
4	Benzene, 2-propenyl-	118	C9H10	000300-57-2	90
5	Benzene, 1-propenyl-	118	C9H10	000637-50-3	90





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060416\
Data File : BF087696.D
Acq On : 5 Jun 2016 8:41
Operator : UM/SJ
Sample : H3345-01
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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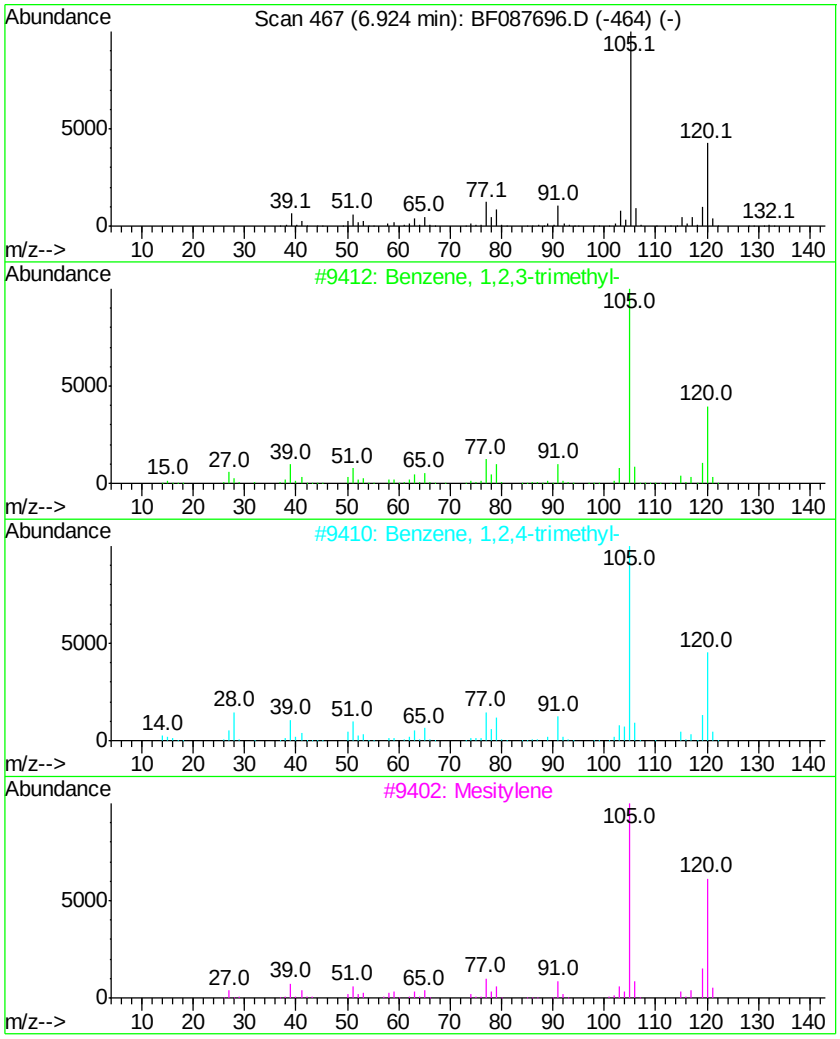
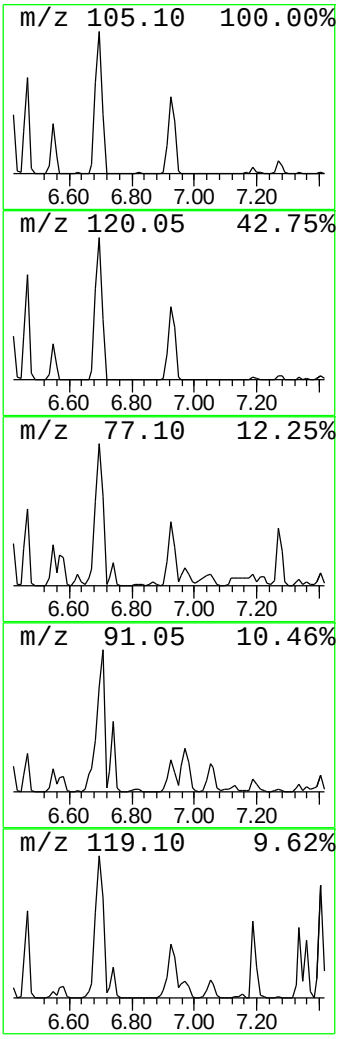
Peak Number 14 Benzene, 1,2,3-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.92	69.28 ng	2653880	1,4-Dichlorobenzene-d4	6.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	93
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	90





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Operator : UM/SJ
Sample : H3345-01
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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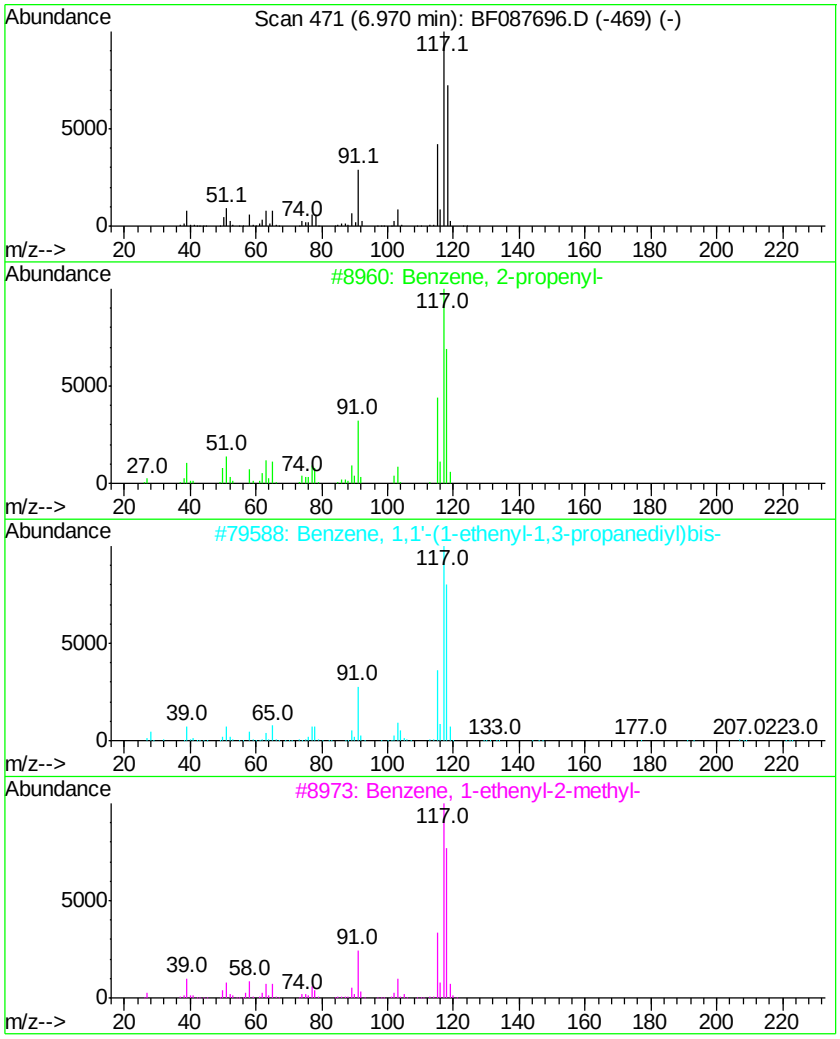
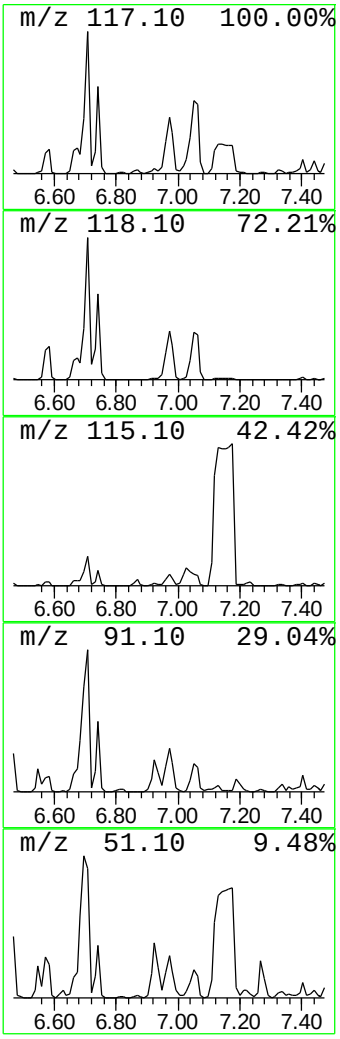
Peak Number 15 Benzene, 2-propenyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.97	52.41 ng	2007390	1,4-Dichlorobenzene-d4	6.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-propenyl-	118	C9H10	000300-57-2	89
2	Benzene, 1,1'-(1-ethenyl-1,3-pro...	222	C17H18	061141-97-7	86
3	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	86
4	Benzene, 1-propenyl-	118	C9H10	000637-50-3	86
5	Indane	118	C9H10	000496-11-7	76





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Acq On : 5 Jun 2016 8:41
Operator : UM/SJ
Sample : H3345-01
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

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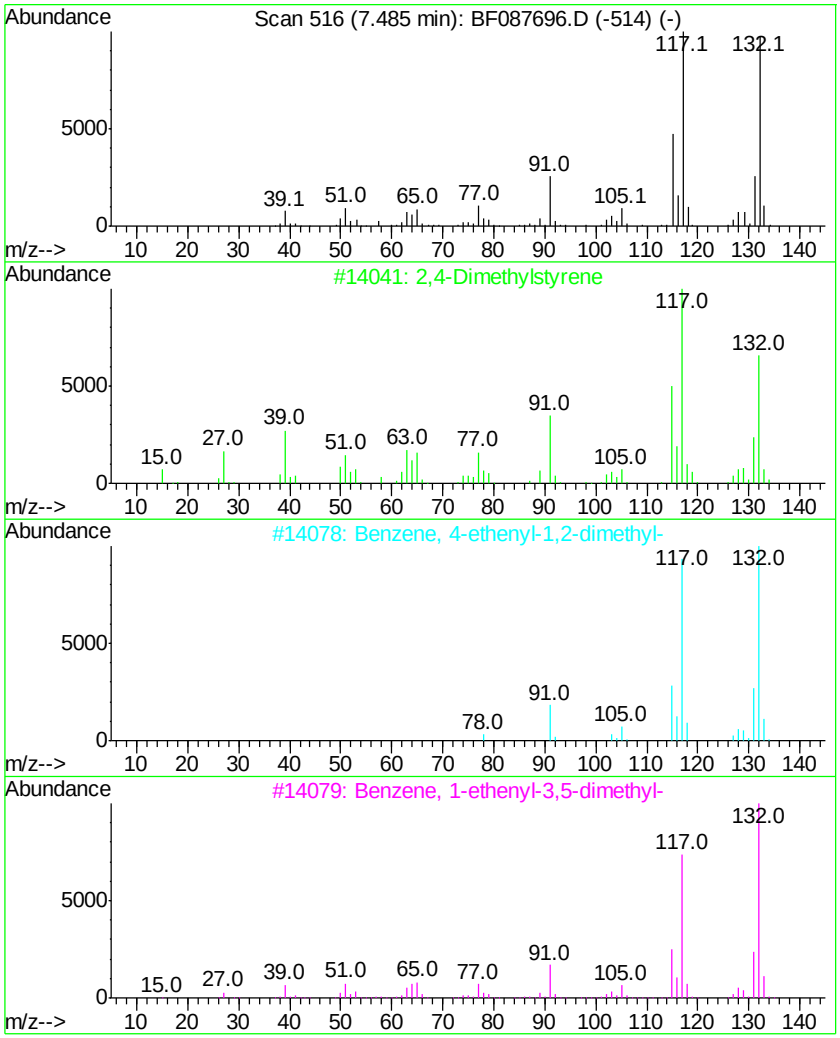
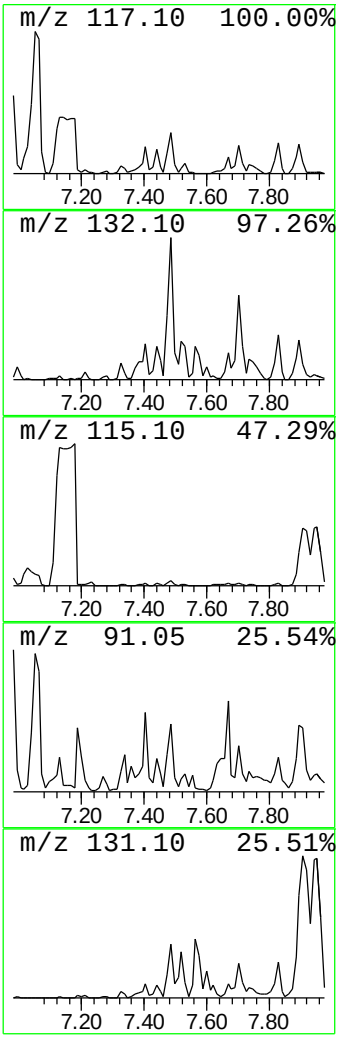
Peak Number 17 2,4-Dimethylstyrene Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.48	16.55 ng	634134	1,4-Dichlorobenzene-d4	6.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4-Dimethylstyrene	132	C10H12	002234-20-0	96
2	Benzene, 4-ethenyl-1,2-dimethyl-	132	C10H12	027831-13-6	96
3	Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	95
4	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	95
5	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	94





Library Search Compound Report

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Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060416.M
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TIC Library : C:\Database\NIST11.L
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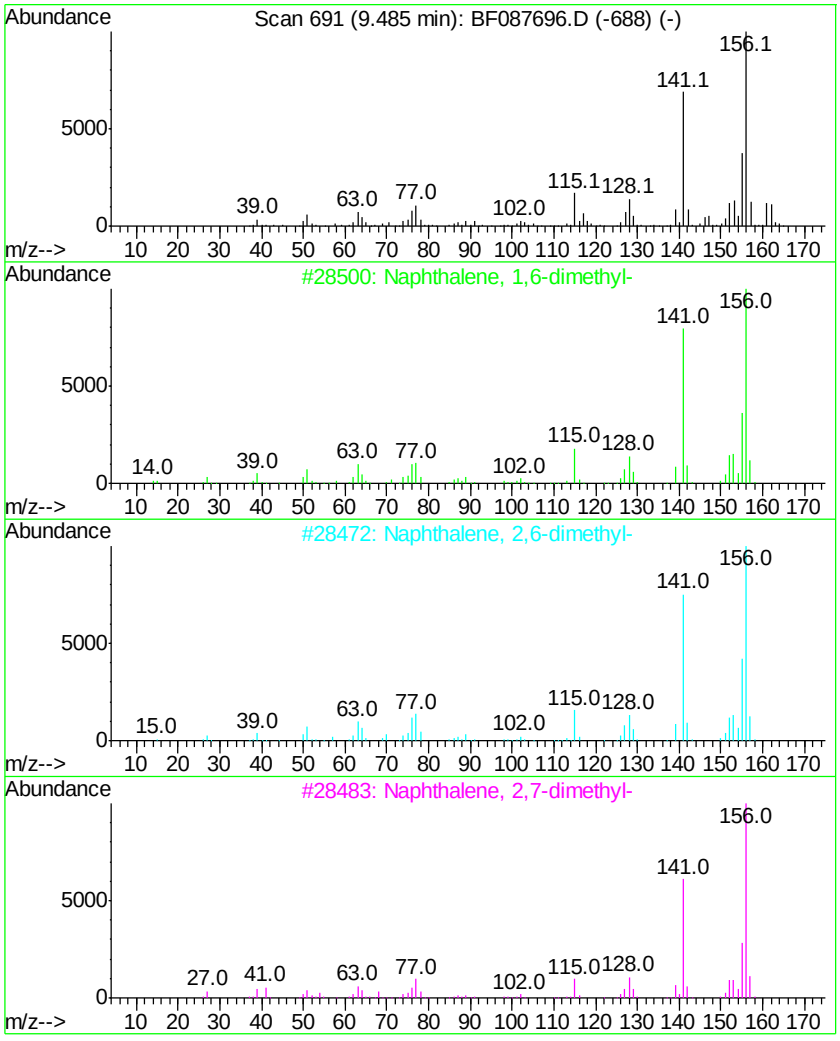
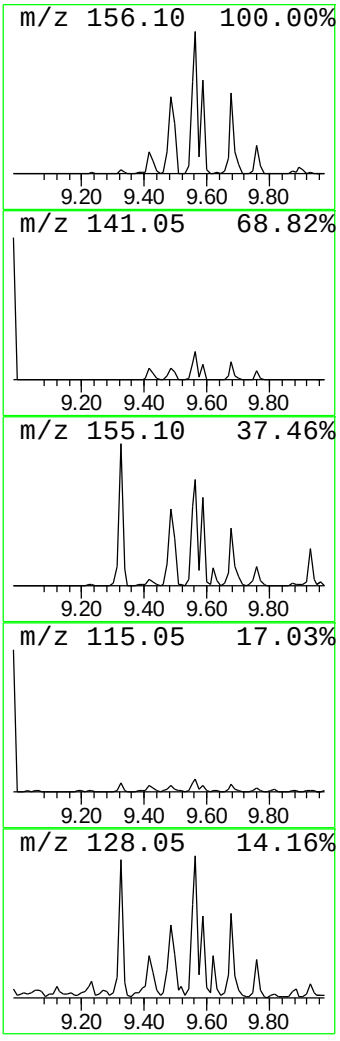
Peak Number 19 Naphthalene, 1,6-dimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.48	14.08 ng	990057	Acenaphthene-d10	9.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
2	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
3	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
4	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
5	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96





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Acq On : 5 Jun 2016 8:41
Operator : UM/SJ
Sample : H3345-01
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

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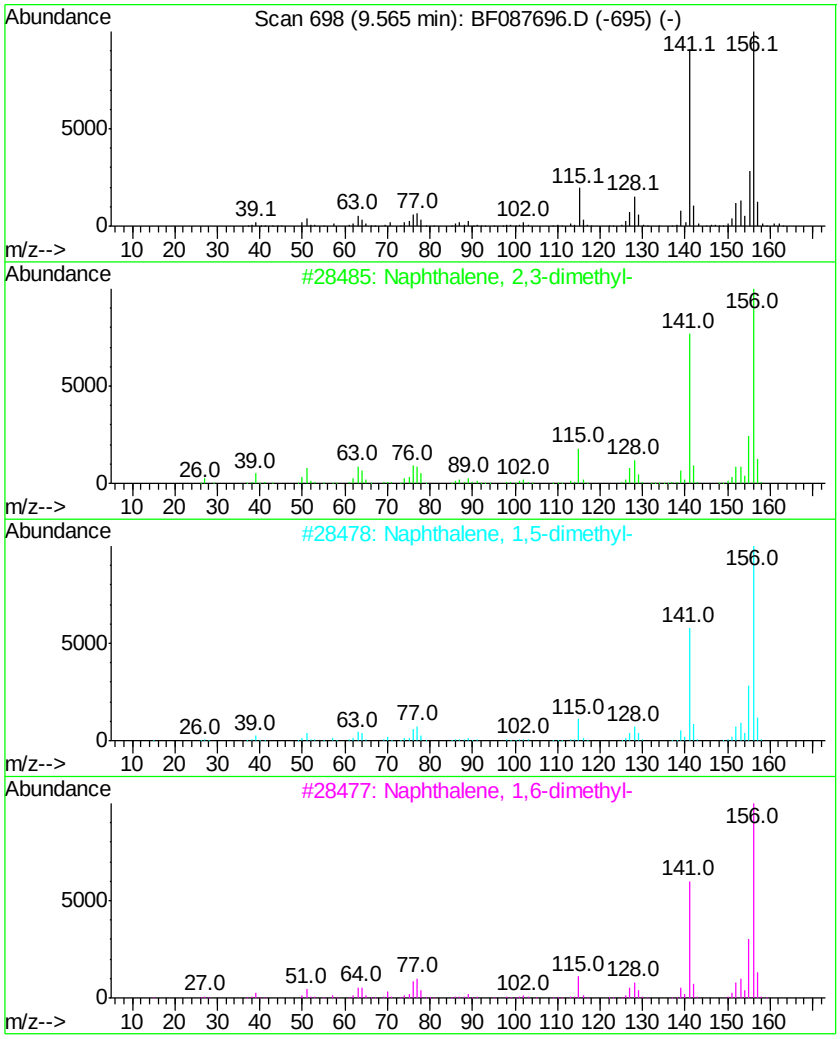
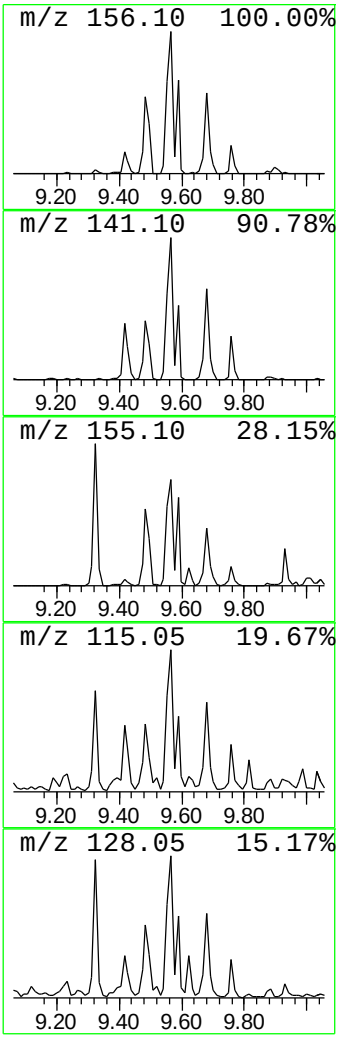
Peak Number 20 Naphthalene, 2,3-dimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.56	21.27 ng	1495350	Acenaphthene-d10	9.90

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
3	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
4	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
5	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060416\
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Operator : UM/SJ
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Misc :
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Instrument : BNA_F
ClientSampleId : MW-16

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Bicyclo[4.2.0]oct...	5.72	448.0	ng	17162300	1	6.87	766100	20.0
Benzene, propyl-	6.32	17.1	ng	656940	1	6.87	766100	20.0
Benzene, 1-ethyl-...	6.39	12.8	ng	489829	1	6.87	766100	20.0
Benzene, 1-ethyl-...	6.42	34.3	ng	1313830	1	6.87	766100	20.0
.alpha.-Methylsty...	6.58	26.9	ng	1029300	1	6.87	766100	20.0
unknown6.63	6.63	77.0	ng	2950150	1	6.87	766100	20.0
Benzene, 1-etheny...	6.74	44.2	ng	1694000	1	6.87	766100	20.0
Benzene, 1,2,3-tr...	6.92	69.3	ng	2653880	1	6.87	766100	20.0
Benzene, 2-propenyl-	6.97	52.4	ng	2007390	1	6.87	766100	20.0
2,4-Dimethylstyrene	7.48	16.6	ng	634134	1	6.87	766100	20.0
Naphthalene, 1,6-...	9.48	14.1	ng	990057	3	9.90	1406380	20.0
Naphthalene, 2,3-...	9.56	21.3	ng	1495350	3	9.90	1406380	20.0