LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040816\
Data File : BF086189.D
Acq On : 9 Apr 2016 1:05
Operator : UM/SJ
Sample : H2334-03
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TW-3-04062016

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-BF040216.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	2.533	31	39	43	rBV	113577	286496	1.99%	0.179%
2	3.470	113	121	122	rBV	194859	431039	3.00%	0.269%
3	3.516	122	125	132	rVB	474076	1006927	7.01%	0.629%
4	3.653	132	137	141	rBV	163038	345129	2.40%	0.215%
5	3.836	147	153	155	rBV	296932	612439	4.26%	0.382%
6	3.893	155	158	161	rVB	120574	214197	1.49%	0.134%
7	4.007	163	168	171	rBV2	124735	284716	1.98%	0.178%
8	4.064	171	173	177	rVB	169512	329259	2.29%	0.206%
9	4.167	181	182	184	rBV	247047	398910	2.78%	0.249%
10	4.224	184	187	192	rVB	1294848	2003035	13.94%	1.251%
11	4.338	192	197	201	rVB	366331	557612	3.88%	0.348%
12	4.579	215	218	219	rBV	143030	203939	1.42%	0.127%
13	4.613	219	221	223	rVB	119729	156786	1.09%	0.098%
14	4.693	223	228	230	rBV	507604	775944	5.40%	0.484%
15	4.750	230	233	235	rBV3	298429	817855	5.69%	0.511%
16	4.796	235	237	240	rVB	1234062	1738479	12.10%	1.085%
17	4.853	240	242	247	rBV2	812647	2122910	14.78%	1.325%
18	4.921	247	248	251	rVB2	220902	325587	2.27%	0.203%
19	5.001	251	255	259	rVB	205208	342950	2.39%	0.214%
20	5.081	259	262	263	rBV	1176250	1653815	11.51%	1.033%
21	5.150	266	268	269	rBV	737039	695365	4.84%	0.434%
22	5.173	269	270	272	rBV2	405317	777527	5.41%	0.485%
23	5.207	272	273	276	rVB	1061684	1151663	8.02%	0.719%
24	5.287	276	280	282	rBV2	3812635	8864996	61.72%	5.535%
25	5.321	282	283	286	rVB	857714	928506	6.46%	0.580%
26	5.379	286	288	291	rBV	697084	954255	6.64%	0.596%
27	5.424	291	292	293	rVB	281536	193134	1.34%	0.121%
28	5.459	293	295	298	rBV	1114924	2110235	14.69%	1.317%
29	5.504	298	299	301	rVB	716205	927358	6.46%	0.579%
30	5.550	301	303	307	rVB2	3754606	6079325	42.32%	3.796%
31	5.630	307	310	315	rVB	3432079	5841546	40.67%	3.647%
32	5.733	315	319	321	rBV	1772419	2854721	19.87%	1.782%
33	5.767	321	322	328	rVB3	452039	1188291	8.27%	0.742%
34	5.870	328	331	334	rBV2	1538835	2996400	20.86%	1.871%

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Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	5.927	334	336	338	rBV2	535084	971403	6.76%	0.606%
36	5.973	338	340	343	rBV2	2527919	5072830	35.32%	3.167%
37	6.030	343	345	346	rBV	633878	609591	4.24%	0.381%
38	6.179	354	358	361	rBV3	1658091	5393490	37.55%	3.367%
39	6.259	361	365	370	rVV2	3557058	13637091	94.94%	8.514%
40	6.339	370	372	375	rVV2	2439782	4681510	32.59%	2.923%
41	6.430	377	380	382	rVB	2113877	3777300	26.30%	2.358%
42	6.510	382	387	390	rBV5	1328352	4236230	29.49%	2.645%
43	6.579	390	393	400	rVB2	3869366	14363891	100.00%	8.968%
44	6.704	400	404	405	rBV2	1345409	2908322	20.25%	1.816%
45	6.773	408	410	411	rBV	1251904	1889423	13.15%	1.180%
46	6.807	411	413	419	rVB	1834896	3801507	26.47%	2.373%
47	6.899	419	421	422	rBV	1507814	2412523	16.80%	1.506%
48	7.036	427	433	434	rBV4	2299083	6724232	46.81%	4.198%
49	7.082	434	437	441	rVV3	1336746	4220322	29.38%	2.635%
50	7.150	441	443	444	rVV	974010	1372688	9.56%	0.857%
51	7.253	448	452	453	rBV	1006466	2619077	18.23%	1.635%
52	7.322	453	458	460	rBV3	1258597	4235570	29.49%	2.644%
53	7.562	477	479	485	rVB3	1626621	5521551	38.44%	3.447%
54	7.722	487	493	496	rBV3	1207190	5881808	40.95%	3.672%
55	12.374	897	900	901	rVB	1382222	2133855	14.86%	1.332%
56	12.728	929	931	933	rVB	1562326	2446945	17.04%	1.528%
57	12.842	939	941	947	rVB2	1136906	2141339	14.91%	1.337%
58	13.082	959	962	963	rBV	1949671	2237004	15.57%	1.397%
59	13.414	989	991	998	rVB	1893048	2313735	16.11%	1.445%
60	13.711	1015	1017	1018	rBV	223613	317745	2.21%	0.198%
61	13.734	1018	1019	1025	rVB	1223927	1651196	11.50%	1.031%
62	13.894	1031	1033	1034	rVB	502993	439976	3.06%	0.275%
63	14.054	1044	1047	1049	rBV	592995	779000	5.42%	0.486%
64	14.351	1071	1073	1075	rVB	433879	422918	2.94%	0.264%
65	14.648	1096	1099	1101	rVB	142187	202255	1.41%	0.126%
66	15.277	1152	1154	1157	rBV	472780	583654	4.06%	0.364%

Sum of corrected areas: 160171327

LSC Report - Integrated Chromatogram

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Operator : UM/SJ
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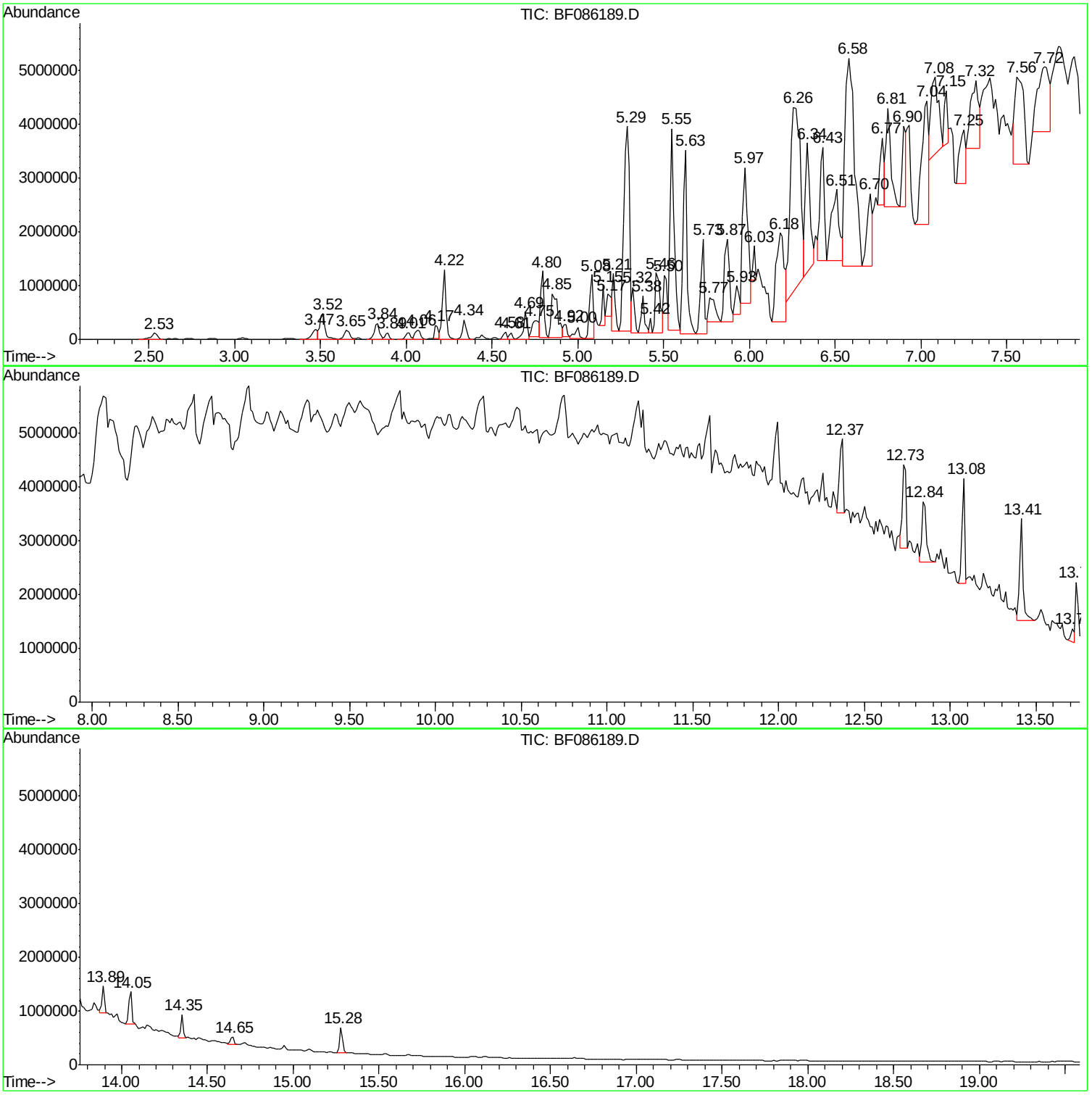
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TW-3-04062016

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.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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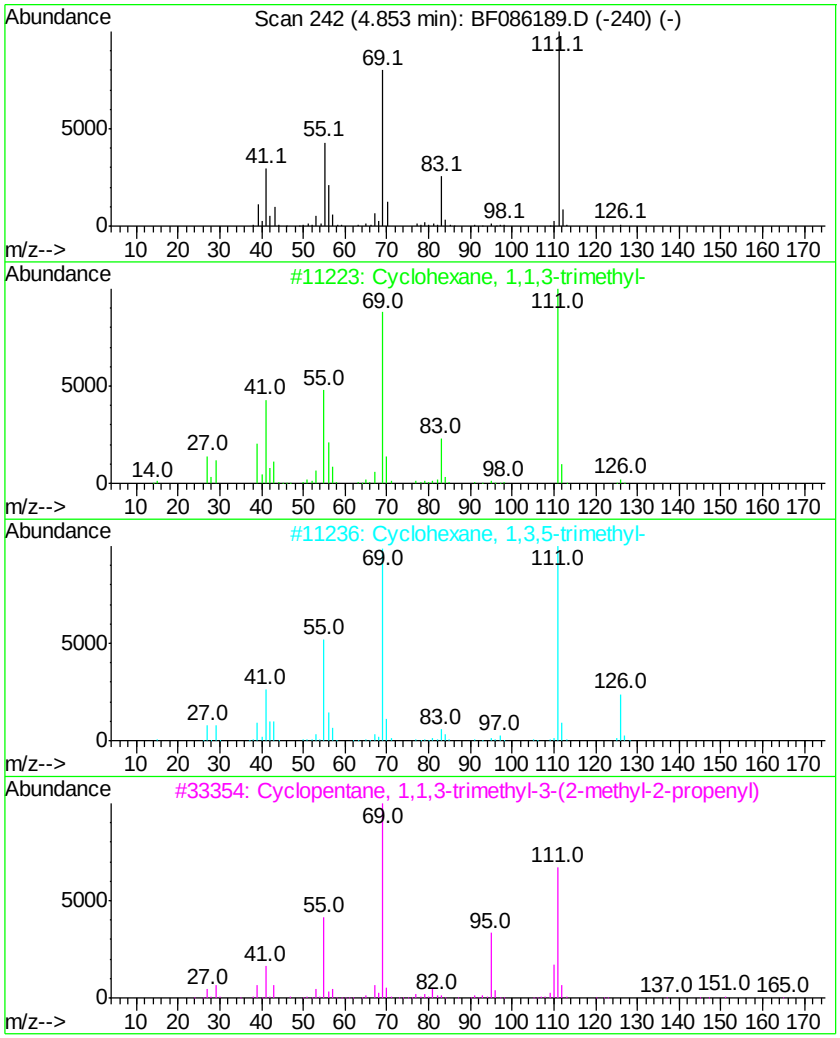
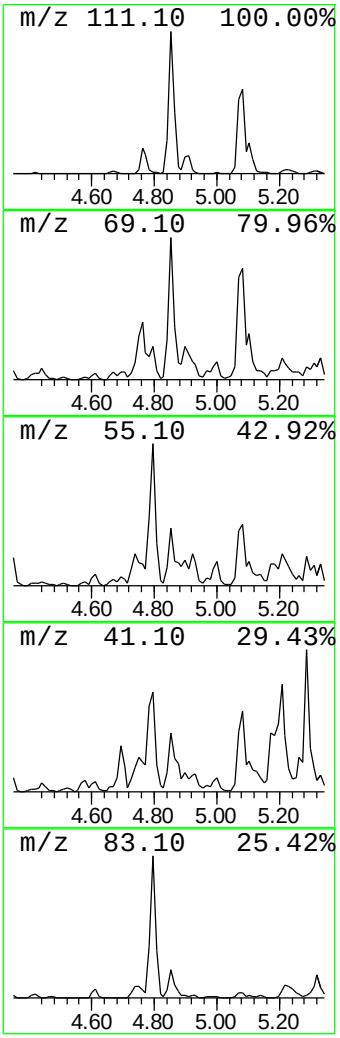
****

Peak Number 1 Cyclohexane, 1,1,3-trimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.85	22.47 ng	2122910	1,4-Dichlorobenzene-d4	6.75

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	90
2	Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	72
3	Cyclopentane, 1,1,3-trimethyl-3-...	166	C12H22	074421-09-3	59
4	Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-27-3	56
5	Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-26-2	56





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040816\
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Acq On : 9 Apr 2016 1:05
Operator : UM/SJ
Sample : H2334-03
Misc :
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Instrument : BNA_F
ClientSampleId : TW-3-04062016

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.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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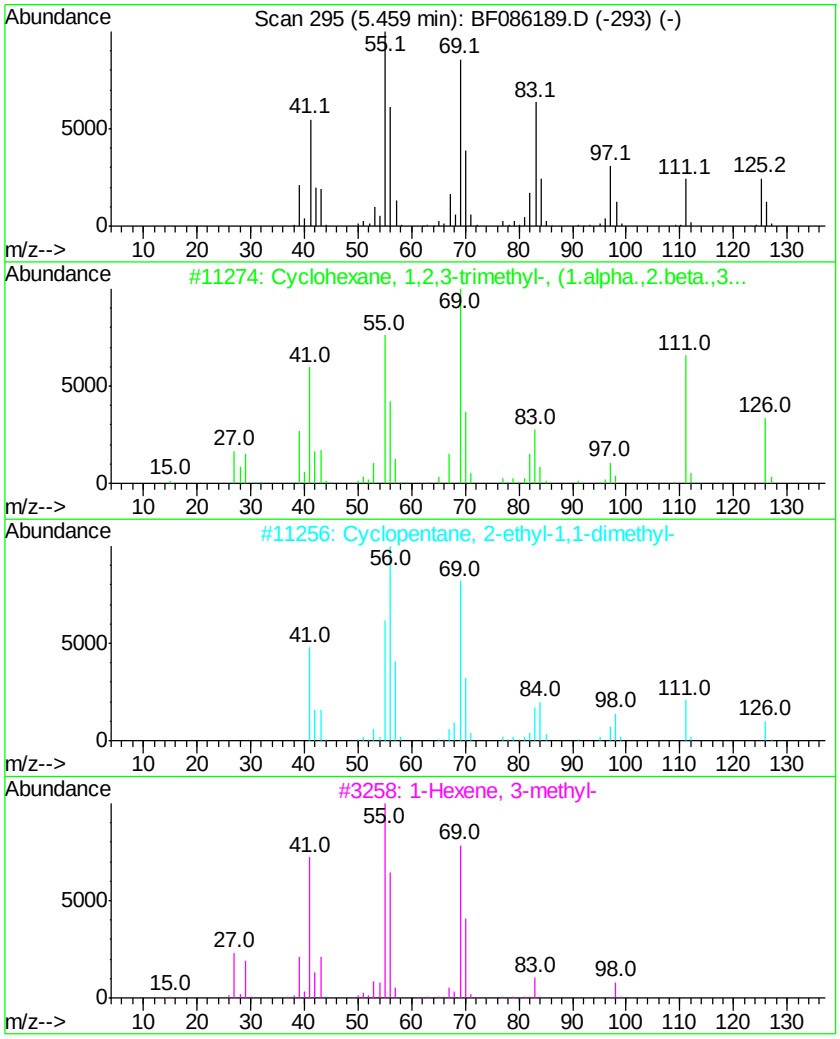
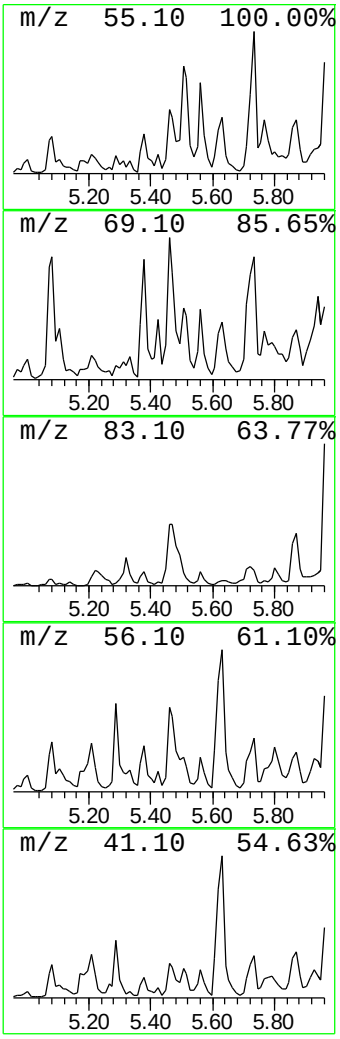
Peak Number 2 Cyclohexane, 1,2,3-trimethy... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.46	22.34 ng	2110240	1,4-Dichlorobenzene-d4	6.75

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,2,3-trimethyl-, (...	126	C9H18	001678-81-5	62
2	Cyclopentane, 2-ethyl-1,1-dimethyl-	126	C9H18	054549-80-3	49
3	1-Hexene, 3-methyl-	98	C7H14	003404-61-3	49
4	Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	46
5	2,2-Dimethyl-3-heptene trans	126	C9H18	019550-75-5	46





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Operator : UM/SJ
Sample : H2334-03
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TW-3-04062016

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.M
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TIC Library : C:\DATABASE\NIST02.L
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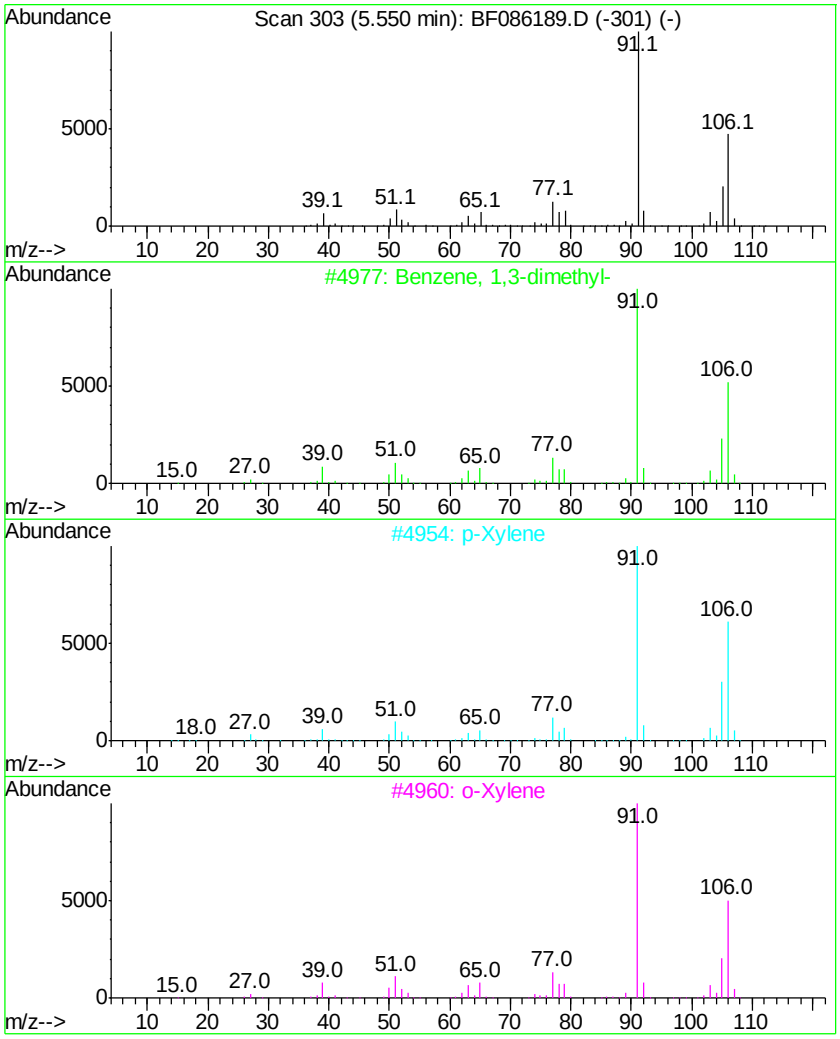
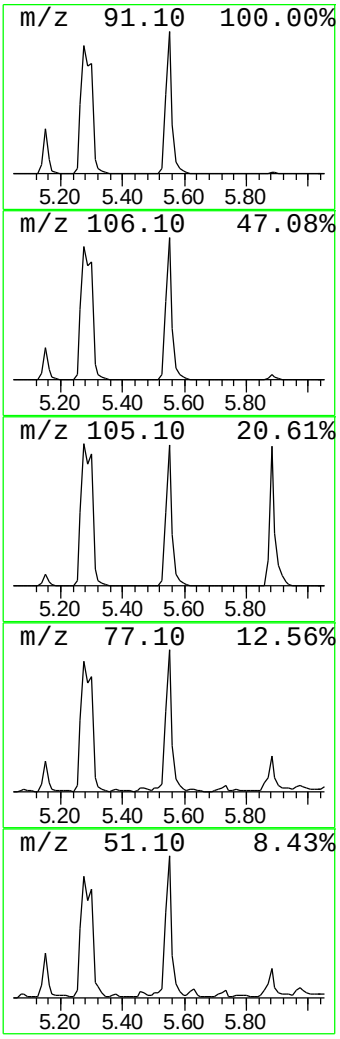
Peak Number 3 Benzene, 1,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.55	64.35 ng	6079330	1,4-Dichlorobenzene-d4	6.75

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2	p-Xylene	106	C8H10	000106-42-3	97
3	o-Xylene	106	C8H10	000095-47-6	97
4	1,3-Cyclopentadiene, 5-(1-methyl...	106	C8H10	002175-91-9	87
5	Ethylbenzene	106	C8H10	000100-41-4	83





Library Search Compound Report

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ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TW-3-04062016

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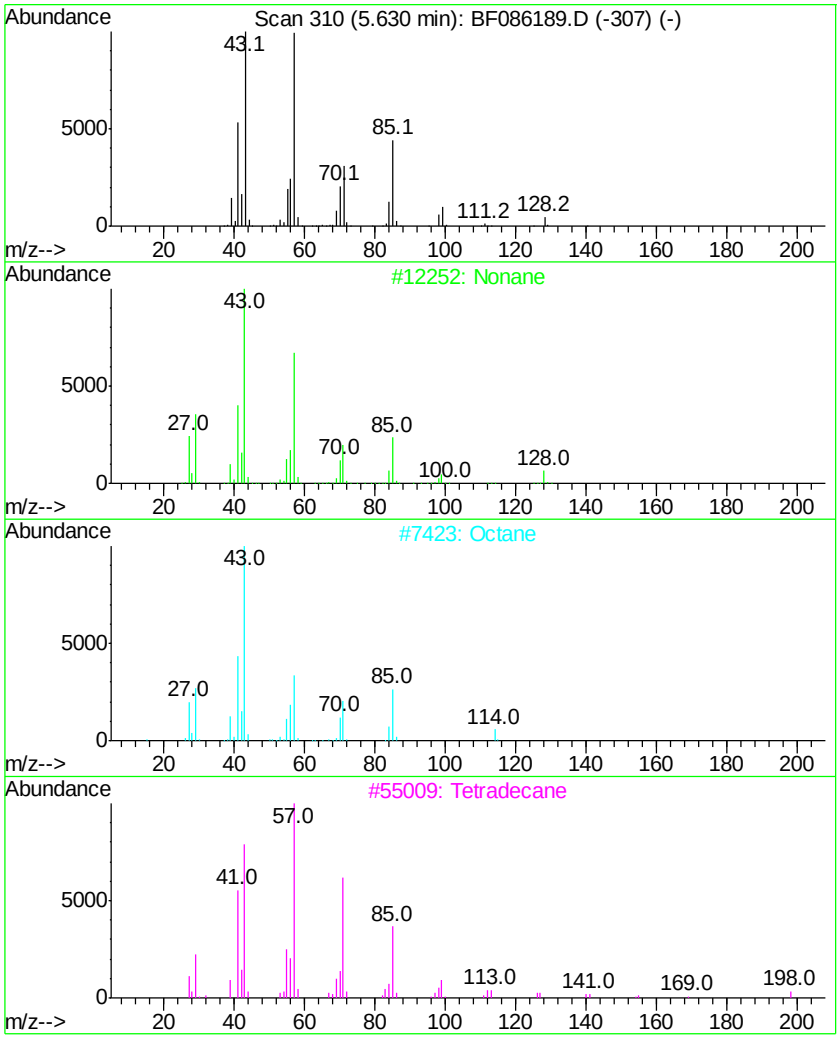
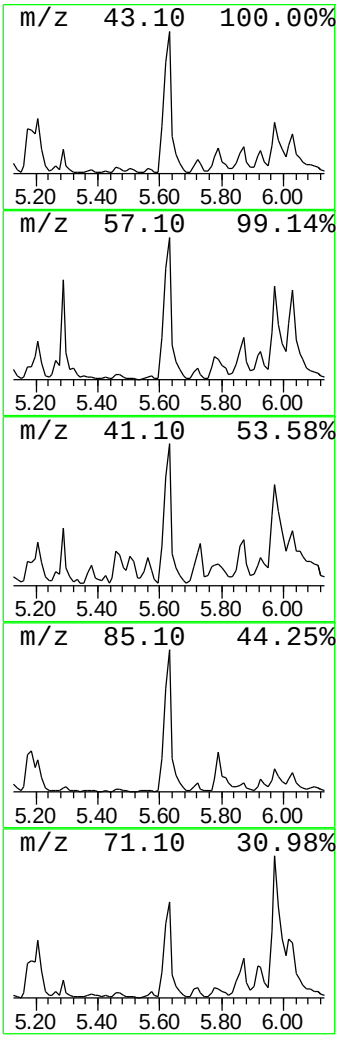
Peak Number 4 Nonane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.63	61.83 ng	5841550	1,4-Dichlorobenzene-d4	6.75

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane	128	C9H20	000111-84-2	94
2	Octane	114	C8H18	000111-65-9	72
3	Tetradecane	198	C14H30	000629-59-4	72
4	4,4-Dimethyl octane	142	C10H22	015869-95-1	59
5	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	47





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Sample : H2334-03
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ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TW-3-04062016

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
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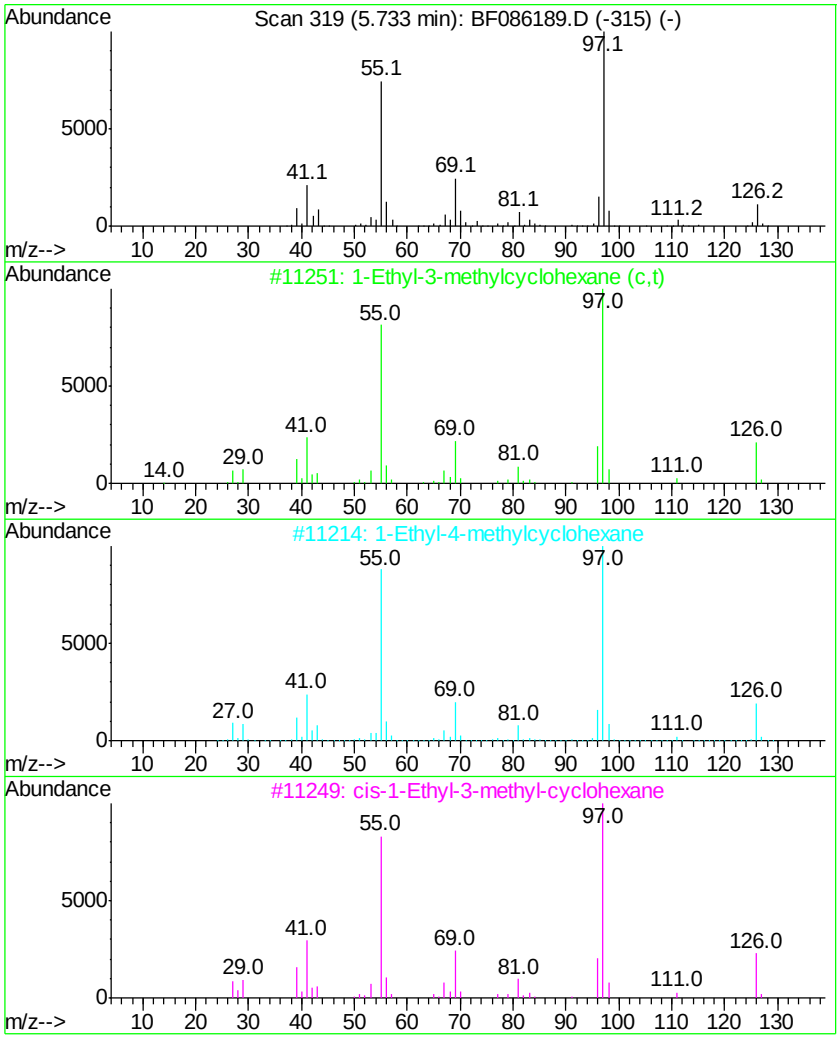
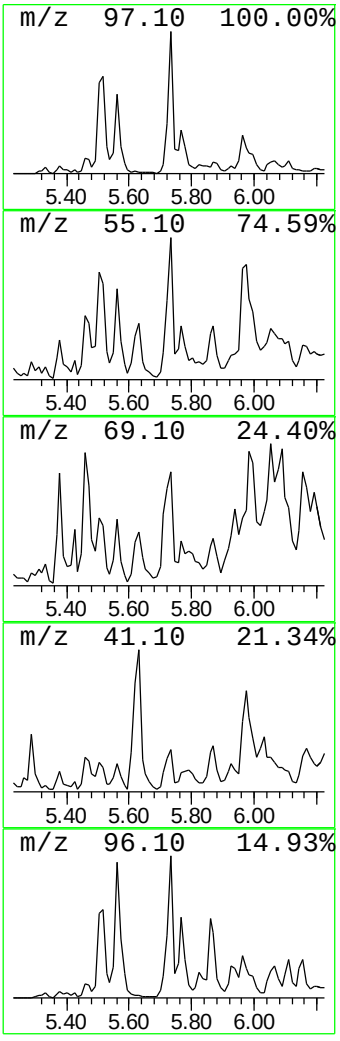
Peak Number 5 1-Ethyl-3-methylcyclohexane... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.73	30.22 ng	2854720	1,4-Dichlorobenzene-d4	6.75

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	94
2	1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	91
3	cis-1-Ethyl-3-methyl-cyclohexane	126	C9H18	019489-10-2	91
4	Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	90
5	Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	80





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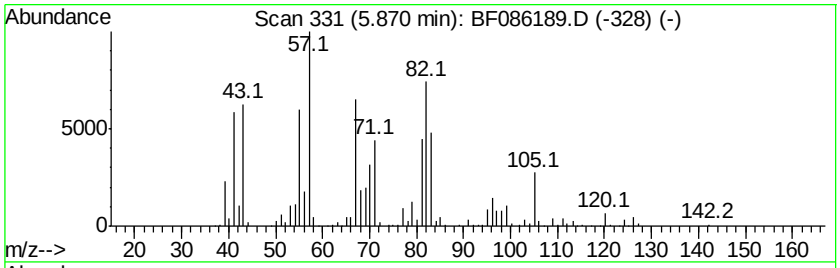
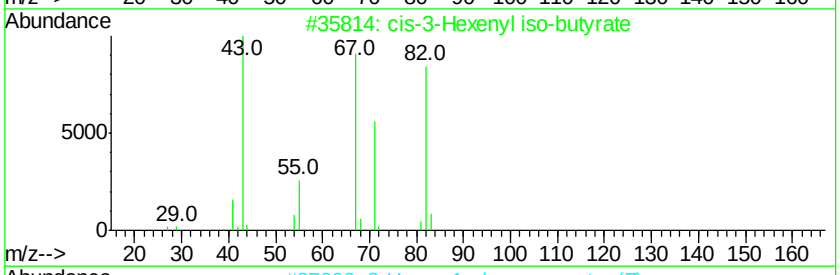
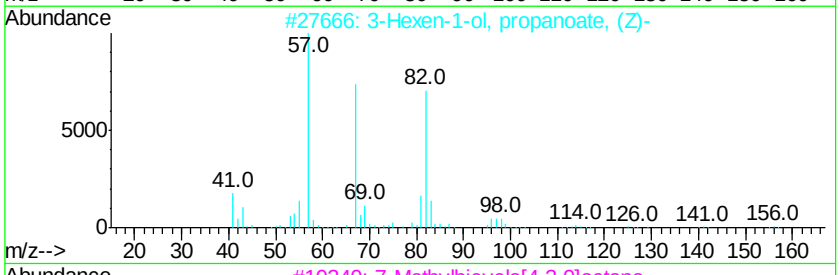
Peak Number 6 unknown5.87 Concentration Rank 15

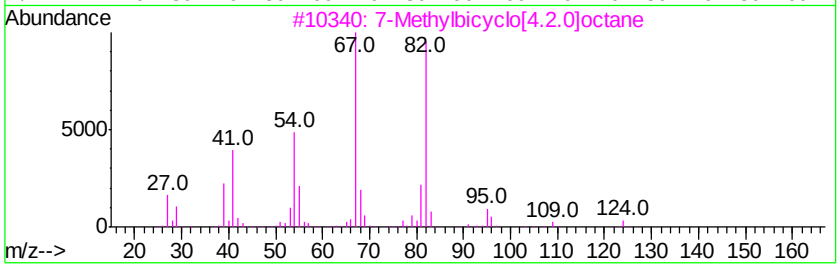
R.T.	EstConc	Area	Relative to ISTD	R.T.
5.87	31.72 ng	2996400	1,4-Dichlorobenzene-d4	6.75

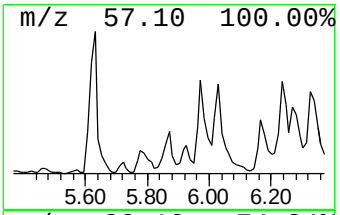
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	cis-3-Hexenyl iso-butyrate	170	C10H18O2	041519-23-7	43
2	3-Hexen-1-ol, propanoate, (Z)-	156	C9H16O2	033467-74-2	43
3	7-Methylbicyclo[4.2.0]octane	124	C9H16	1000210-90-2	35
4	2,4-Hexadiene, (E,Z)-	82	C6H10	005194-50-3	30
5	3-Hexen-1-ol, acetate, (E)-	142	C8H14O2	003681-82-1	27

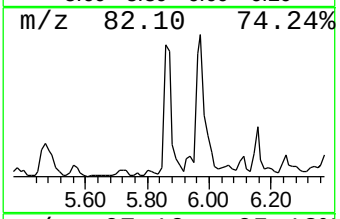


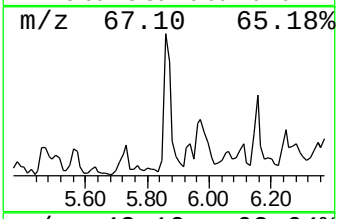


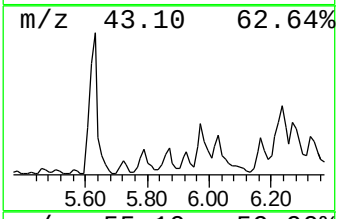


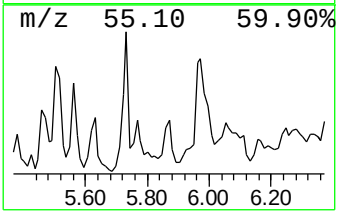












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Operator : UM/SJ
Sample : H2334-03
Misc :
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Instrument : BNA_F
ClientSampleId : TW-3-04062016

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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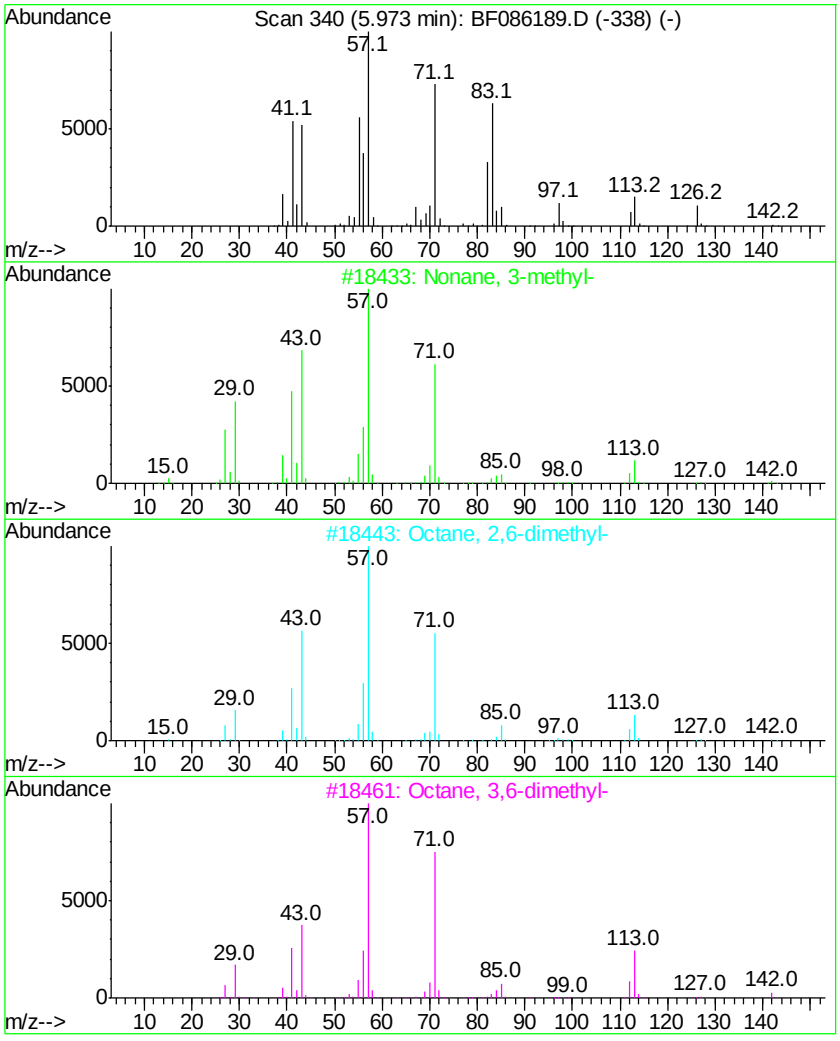
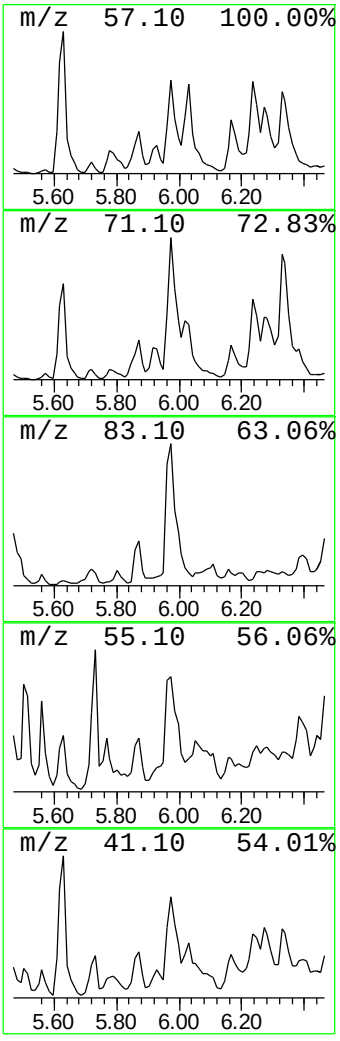
Peak Number 7 Nonane, 3-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.97	53.70 ng	5072830	1,4-Dichlorobenzene-d4	6.75

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane, 3-methyl-	142	C10H22	005911-04-6	55
2	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	55
3	Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	46
4	Cyclohexane, propyl-	126	C9H18	001678-92-8	42
5	3-Bromooctane	192	C8H17Br	000999-64-4	38





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ClientSampleId : TW-3-04062016

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

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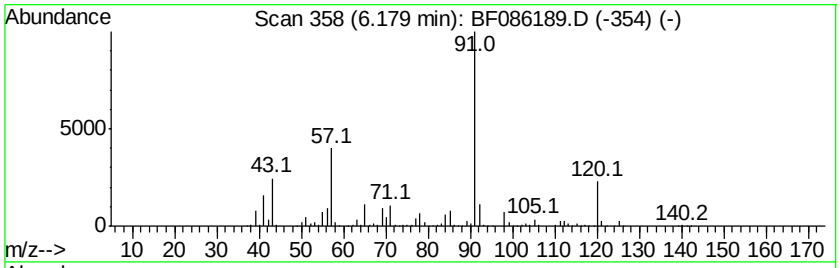
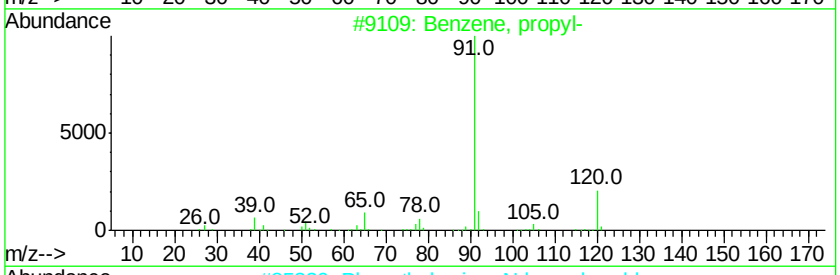
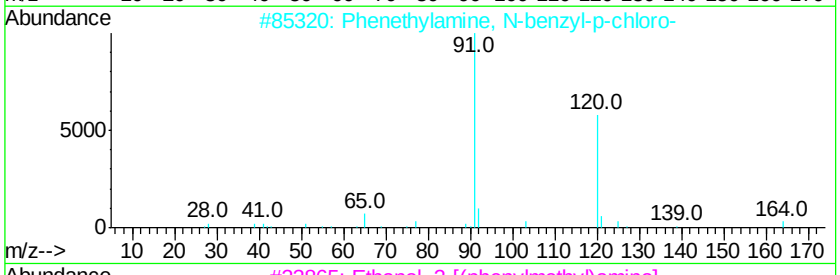
Peak Number 8 Benzene, propyl- Concentration Rank 9

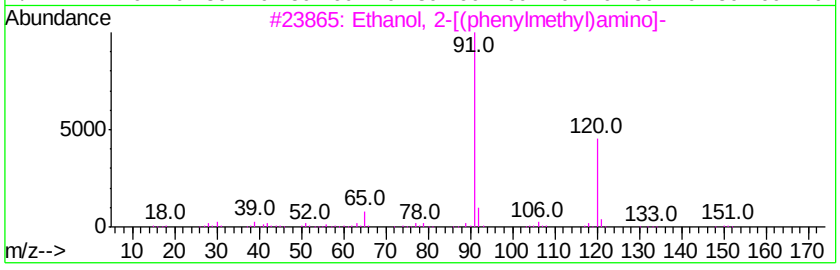
R.T.	EstConc	Area	Relative to ISTD	R.T.
6.18	57.09 ng	5393490	1,4-Dichlorobenzene-d4	6.75

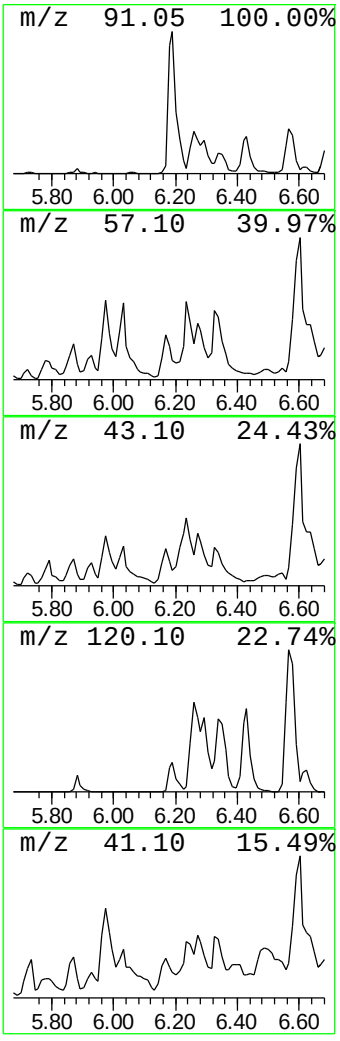
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	55
2	Phenethylamine, N-benzyl-p-chloro-	245	C15H16ClN	013622-43-0	47
3	Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	43
4	N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	43
5	1,2-Ethanediamine, N-(phenylmeth...	150	C9H14N2	004152-09-4	43











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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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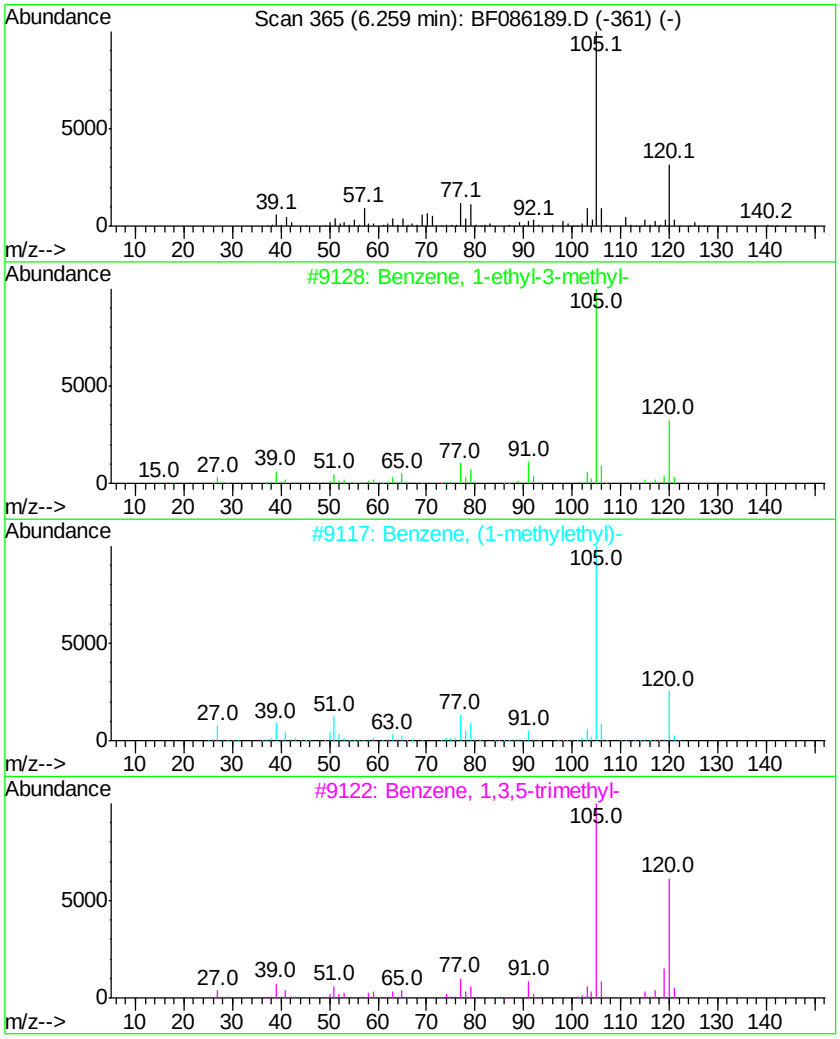
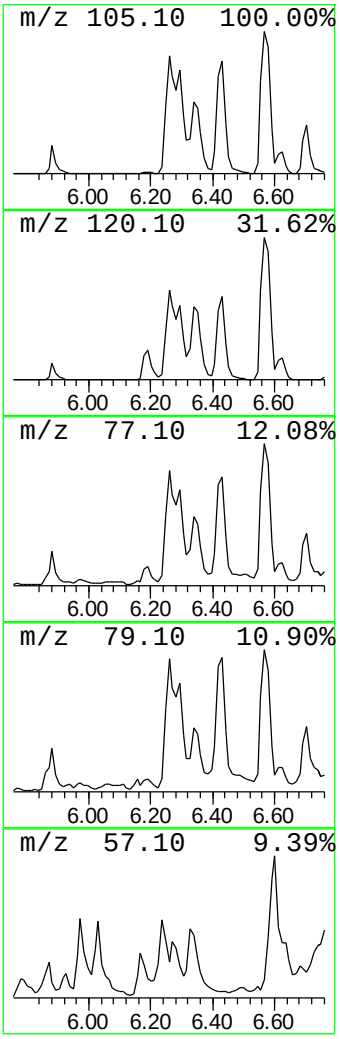
Peak Number 9 Benzene, 1-ethyl-3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.26	144.35 ng	13637100	1,4-Dichlorobenzene-d4	6.75

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
2	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	81
3	Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	81
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	81
5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	81





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

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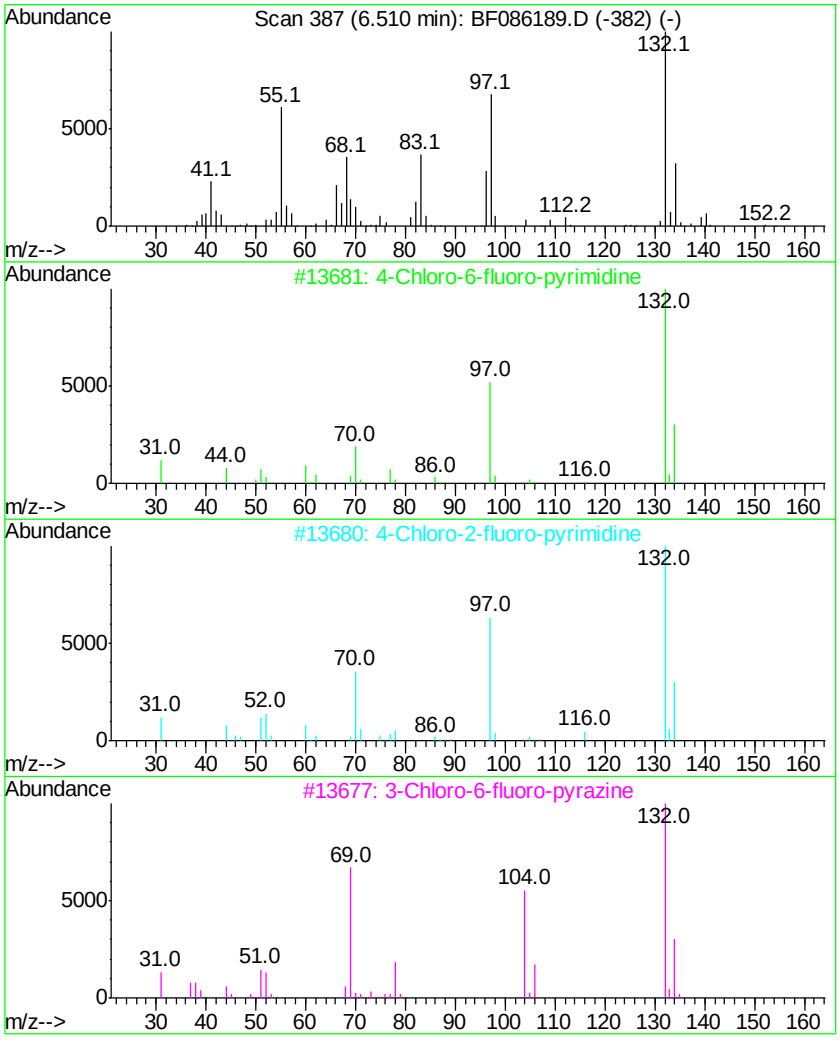
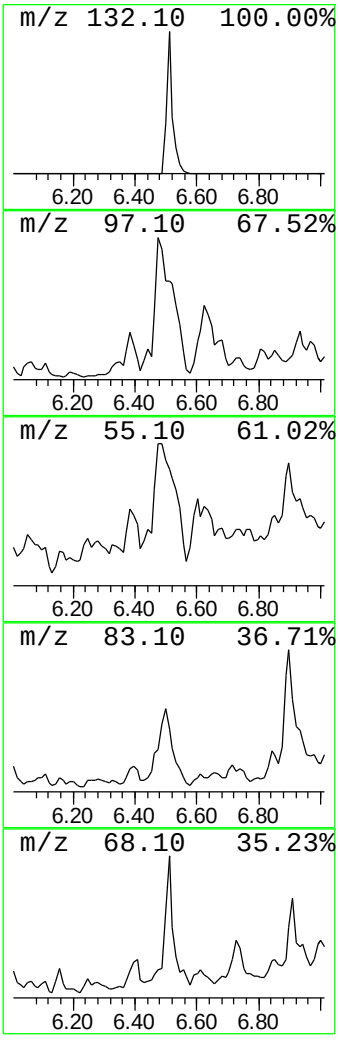
Peak Number 10 unknown6.51 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.51	44.84 ng	4236230	1,4-Dichlorobenzene-d4	6.75

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	25
2	4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	25
3	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
4	Sorbic Acid	112	C6H8O2	000110-44-1	16
5	Cyclohexane, 1-methyl-2-propyl-	140	C10H20	004291-79-6	15





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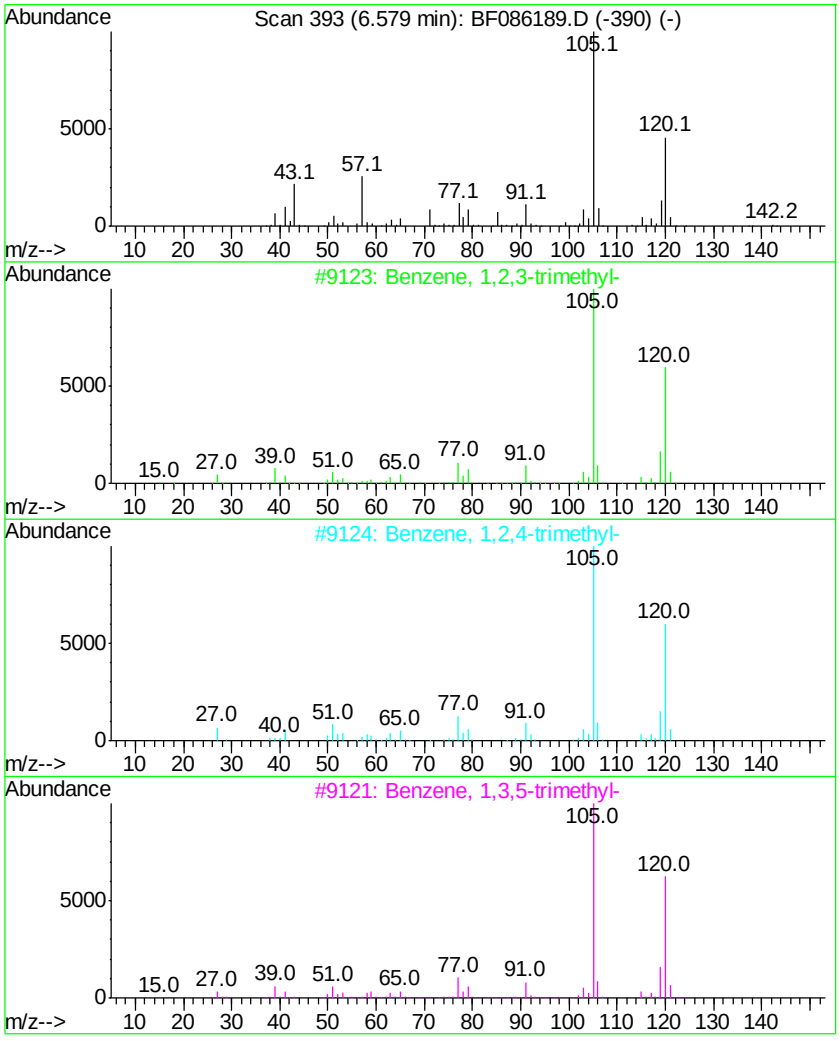
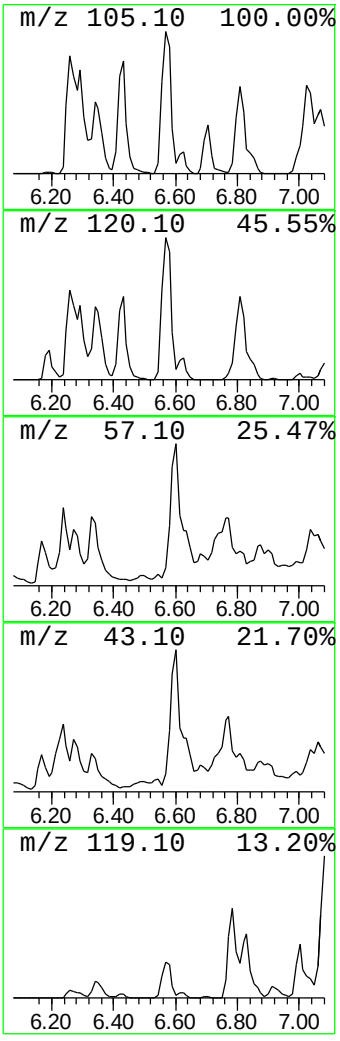
Peak Number 11 Benzene, 1,2,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	152.05 ng	14363900	1,4-Dichlorobenzene-d4	6.75

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
3	Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90





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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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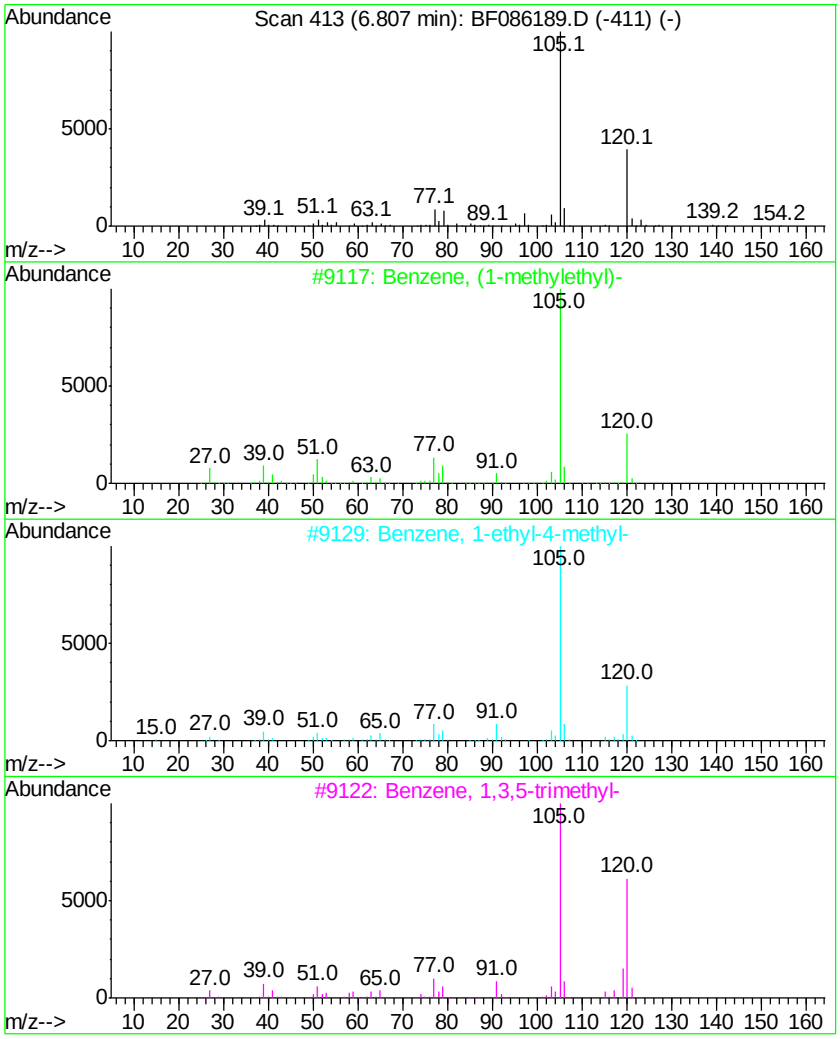
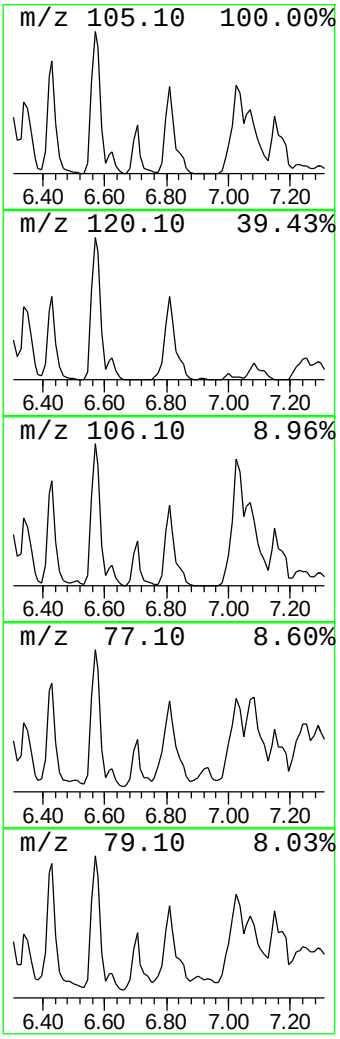
Peak Number 12 Benzene, (1-methylethyl)- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.81	40.24 ng	3801510	1,4-Dichlorobenzene-d4	6.75

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	80
2	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	72
3	Benzene, 1,3,5-trimethyl-	120	C9H12	000108-67-8	72
4	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	72
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	64





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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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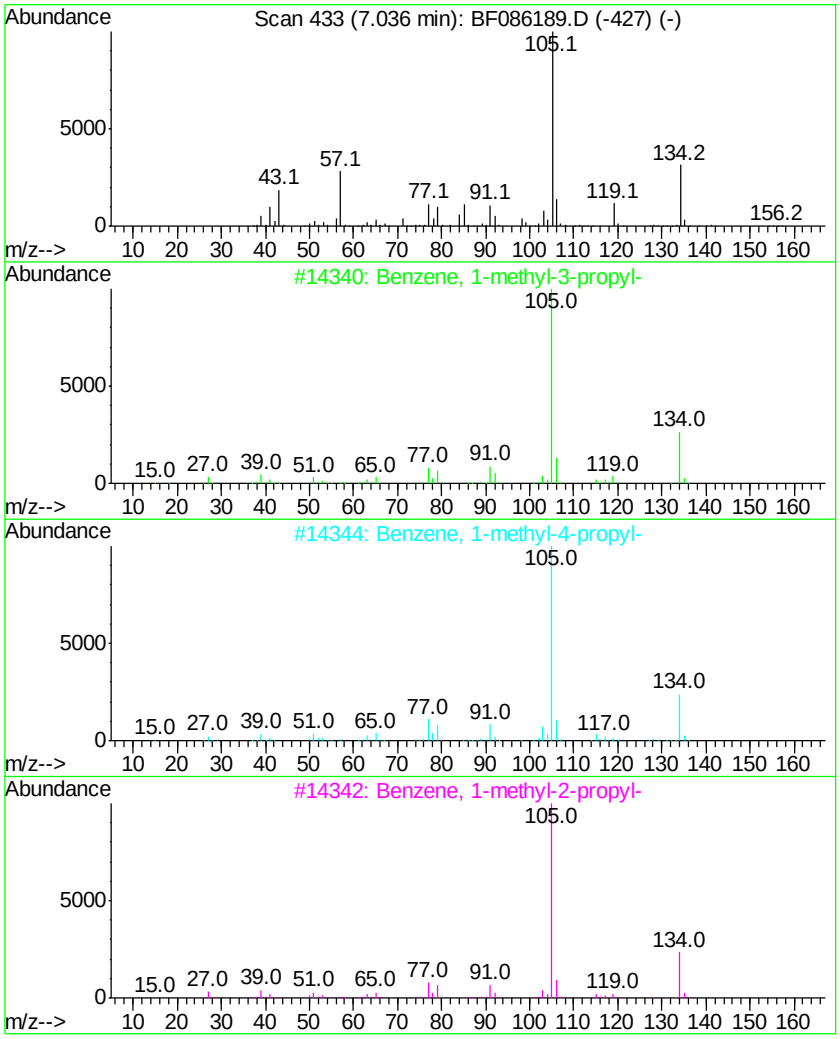
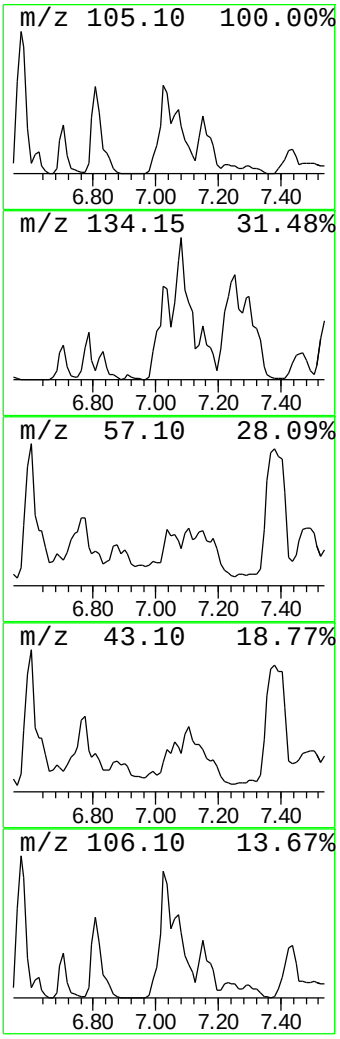
Peak Number 14 Benzene, 1-methyl-3-propyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.04	71.18 ng	6724230	1,4-Dichlorobenzene-d4	6.75

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	76
2	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	76
3	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	76
4	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	58
5	Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	58





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

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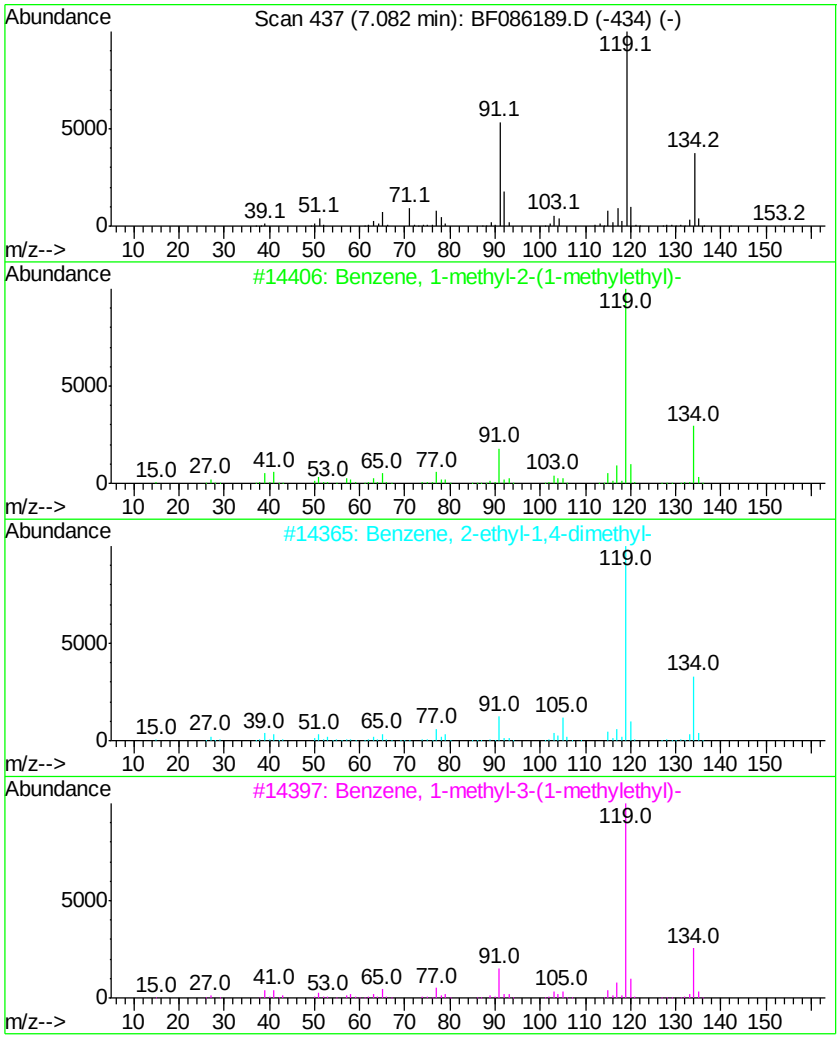
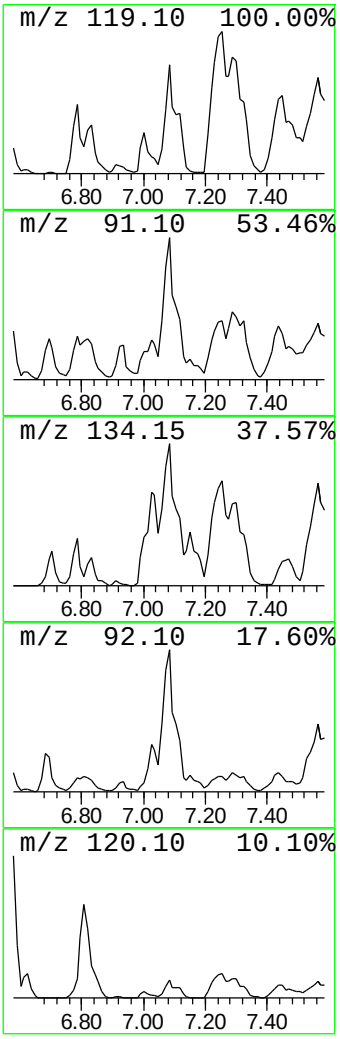
Peak Number 15 Benzene, 1-methyl-2-(1-meth... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.08	44.67 ng	4220320	1,4-Dichlorobenzene-d4	6.75

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	93
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	90
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	90
4	1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	90
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	90





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

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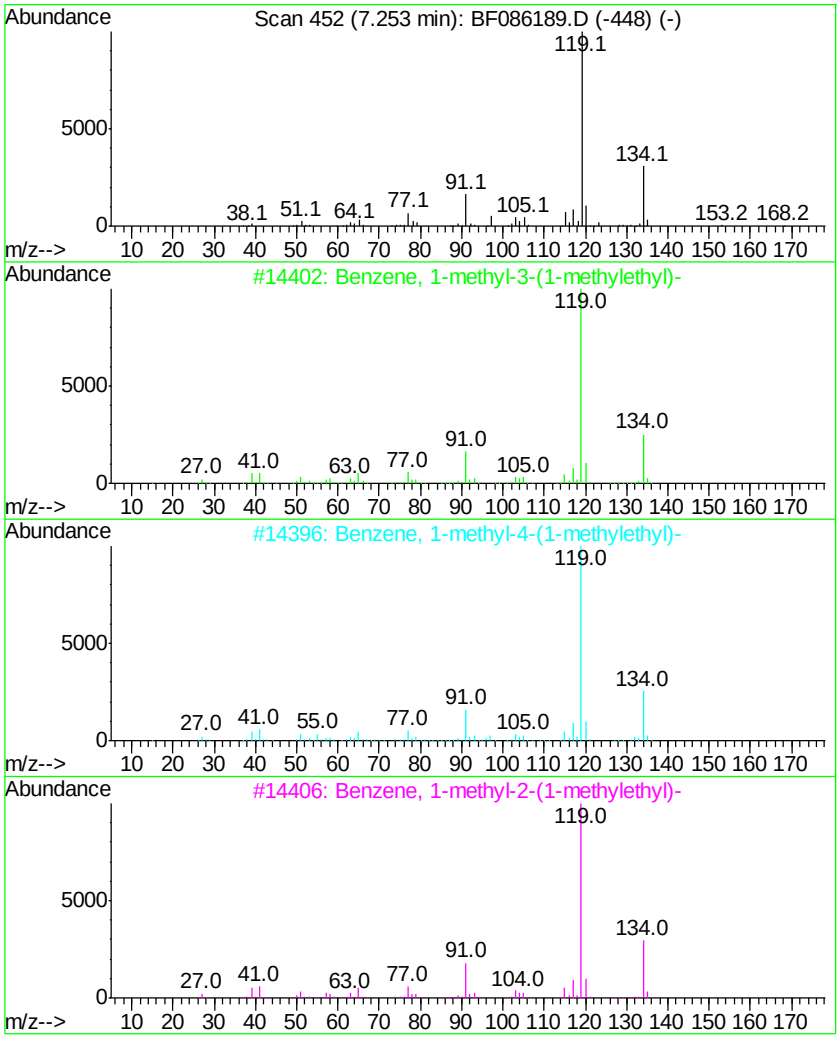
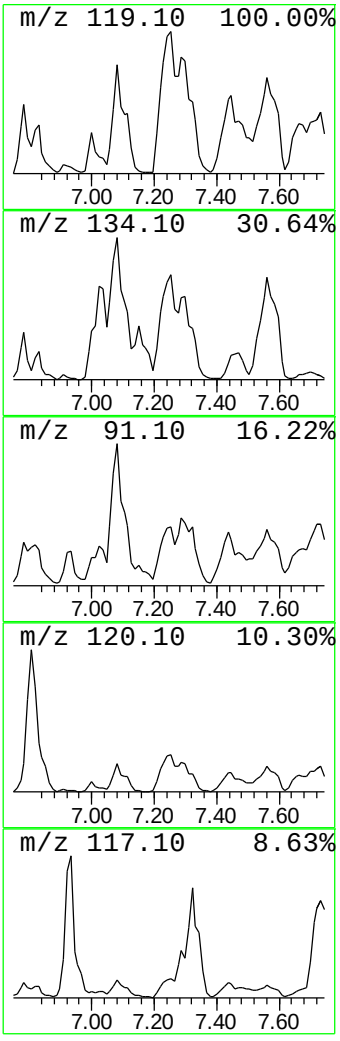
Peak Number 16 Benzene, 1-methyl-3-(1-meth... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.25	27.72 ng	2619080	1,4-Dichlorobenzene-d4	6.75

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
2	Benzene, 1-methyl-4-(1-methyleth...	134	C10H14	000099-87-6	94
3	Benzene, 1-methyl-2-(1-methyleth...	134	C10H14	000527-84-4	93
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	91
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91





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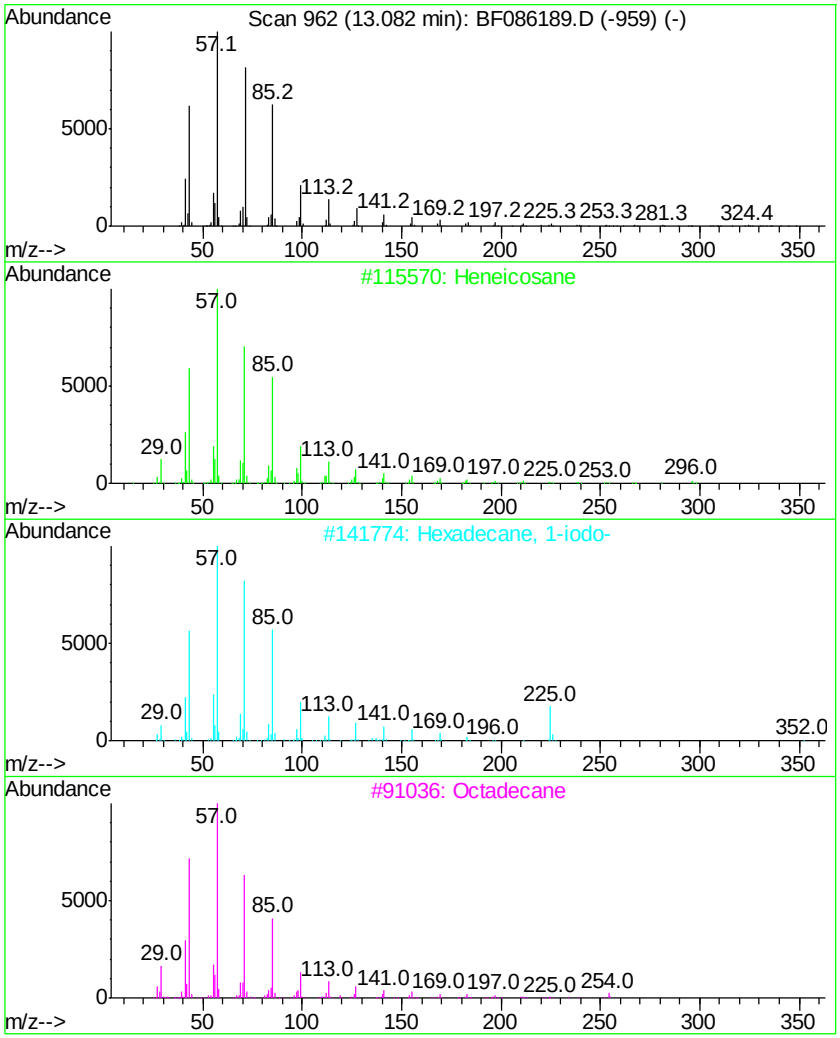
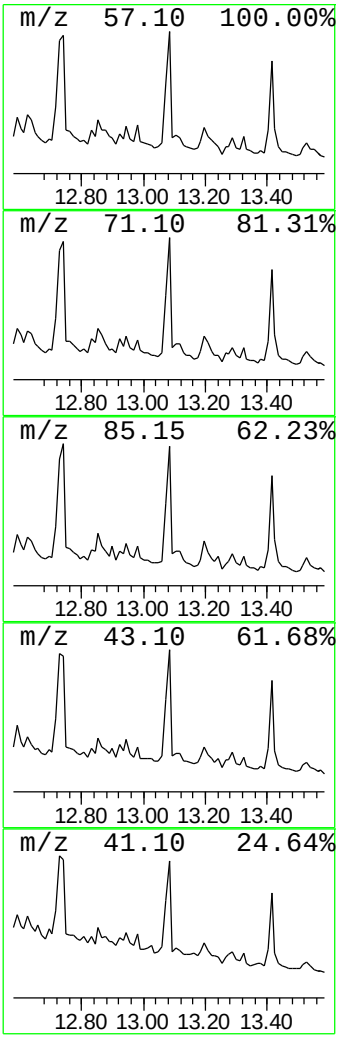
Peak Number 17 Heneicosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.08	101.69 ng	2237000	Chrysene-d12	13.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	90
3	Octadecane	254	C18H38	000593-45-3	86
4	Hexadecane	226	C16H34	000544-76-3	80
5	Tetracosane	338	C24H50	000646-31-1	64





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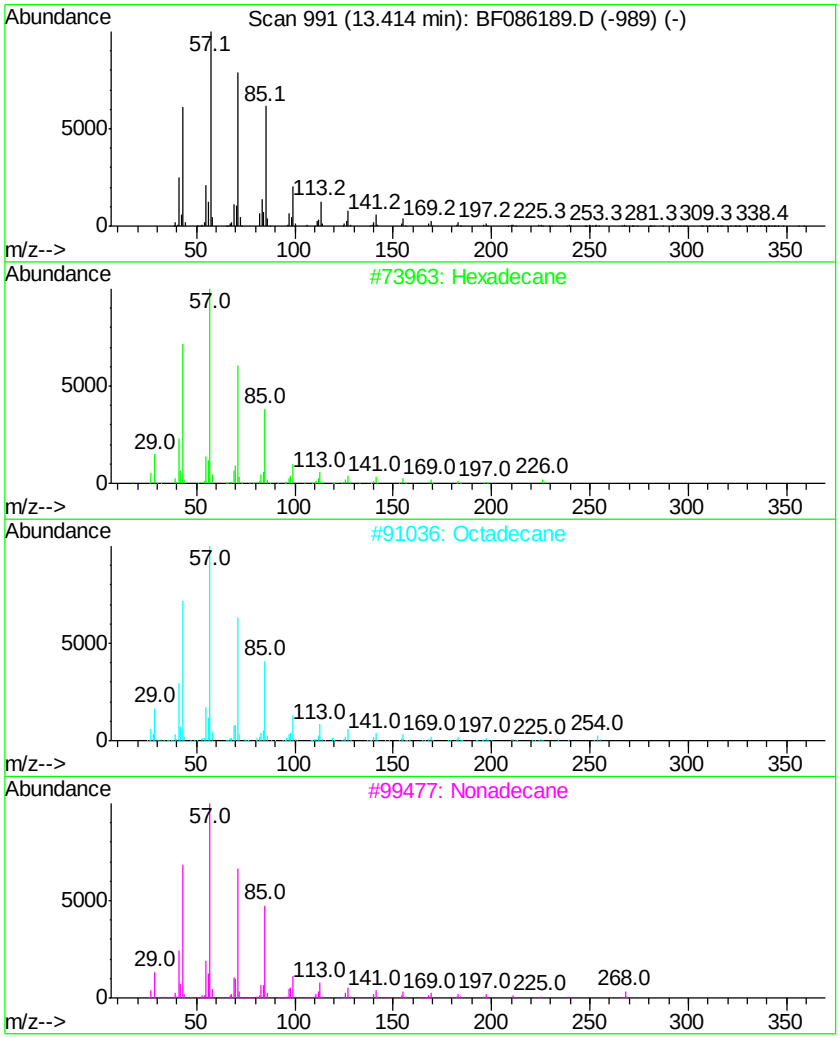
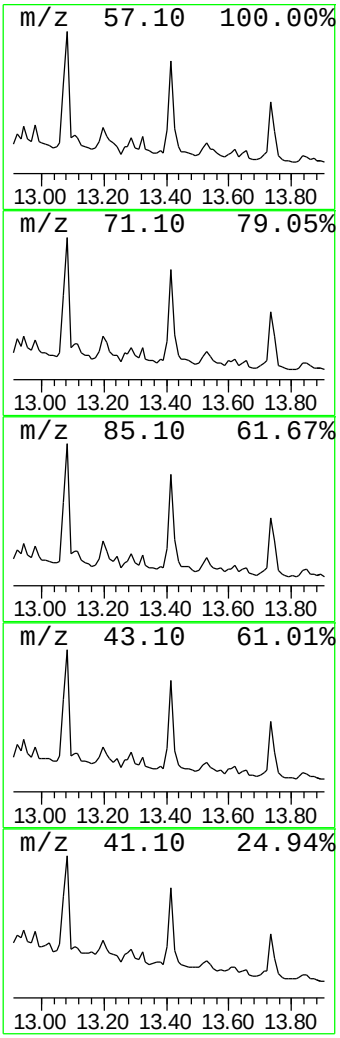
Peak Number 18 Hexadecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.41	105.18 ng	2313740	Chrysene-d12	13.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	90
2	Octadecane	254	C18H38	000593-45-3	86
3	Nonadecane	268	C19H40	000629-92-5	80
4	Dodecane, 1-iodo-	296	C12H25I	004292-19-7	80
5	2-Bromo dodecane	248	C12H25Br	013187-99-0	76





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF040816\
Data File : BF086189.D
Acq On : 9 Apr 2016 1:05
Operator : UM/SJ
Sample : H2334-03
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TW-3-04062016

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040216.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
Cyclohexane, 1,1,...	4.85	22.5	ng	2122910	1	6.75	1889420	20.0
Cyclohexane, 1,2,...	5.46	22.3	ng	2110240	1	6.75	1889420	20.0
Benzene, 1,3-dime...	5.55	64.3	ng	6079330	1	6.75	1889420	20.0
Nonane	5.63	61.8	ng	5841550	1	6.75	1889420	20.0
1-Ethyl-3-methylc...	5.73	30.2	ng	2854720	1	6.75	1889420	20.0
unknown5.87	5.87	31.7	ng	2996400	1	6.75	1889420	20.0
Nonane, 3-methyl-	5.97	53.7	ng	5072830	1	6.75	1889420	20.0
Benzene, propyl-	6.18	57.1	ng	5393490	1	6.75	1889420	20.0
Benzene, 1-ethyl-...	6.26	144.3	ng	13637100	1	6.75	1889420	20.0
unknown6.51	6.51	44.8	ng	4236230	1	6.75	1889420	20.0
Benzene, 1,2,3-tr...	6.58	152.1	ng	14363900	1	6.75	1889420	20.0
Benzene, (1-methy...	6.81	40.2	ng	3801510	1	6.75	1889420	20.0
Benzene, 1-methyl...	7.04	71.2	ng	6724230	1	6.75	1889420	20.0
Benzene, 1-methyl...	7.08	44.7	ng	4220320	1	6.75	1889420	20.0
Benzene, 1-methyl...	7.25	27.7	ng	2619080	1	6.75	1889420	20.0
Heneicosane	13.08	101.7	ng	2237000	5	13.89	439976	20.0
Hexadecane	13.41	105.2	ng	2313740	5	13.89	439976	20.0