LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051416\
Data File : BF087068.D
Acq On : 15 May 2016 14:30
Operator : UM/SJ
Sample : H2968-07 2X
Misc :
ALS Vial : 87 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DUPB

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-BF050916.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	1.735	12	13	52	rVB	2524901	5398484	100.00%	21.636%
2	4.741	273	276	286	rBV	42242	91290	1.69%	0.366%
3	5.198	314	316	329	rBV	1087471	1432391	26.53%	5.741%
4	6.284	408	411	418	rBV	1150668	1413533	26.18%	5.665%
5	6.387	418	420	433	rVB	1470269	1511675	28.00%	6.059%
6	6.616	437	440	442	rBV	331623	440253	8.16%	1.764%
7	6.764	451	453	456	rBV	1249449	1123534	20.81%	4.503%
8	7.187	488	490	492	rBV	827652	844184	15.64%	3.383%
9	7.896	550	552	555	rBV	498927	555344	10.29%	2.226%
10	8.982	644	647	649	rBV	1265471	1375187	25.47%	5.512%
11	9.382	680	682	685	rVB	112852	104805	1.94%	0.420%
12	9.507	692	693	696	rBV	62699	61486	1.14%	0.246%
13	9.587	699	700	703	rBV	49910	65408	1.21%	0.262%
14	9.645	703	705	708	rBV	565133	543151	10.06%	2.177%
15	10.090	742	744	746	rVB	93300	74317	1.38%	0.298%
16	10.433	772	774	777	rBV	688026	706363	13.08%	2.831%
17	10.559	783	785	789	rVB	87244	101599	1.88%	0.407%
18	10.742	798	801	806	rBV2	22616	54391	1.01%	0.218%
19	11.119	832	834	839	rBV2	337737	694073	12.86%	2.782%
20	11.199	839	841	843	rVB	120140	111807	2.07%	0.448%
21	11.622	876	878	880	rBV	57317	82154	1.52%	0.329%
22	11.656	880	881	883	rVV	73153	77069	1.43%	0.309%
23	11.702	883	885	886	rVV	61457	57347	1.06%	0.230%
24	11.736	886	888	893	rVB2	156002	206527	3.83%	0.828%
25	11.919	900	904	908	rBV	53716	98891	1.83%	0.396%
26	12.171	924	926	928	rBV	71594	89121	1.65%	0.357%
27	12.205	928	929	932	rVB3	35586	66456	1.23%	0.266%
28	12.262	932	934	938	rVB2	47661	71573	1.33%	0.287%
29	12.331	938	940	943	rBV	1023532	928721	17.20%	3.722%
30	12.433	947	949	950	rBV	52391	72722	1.35%	0.291%
31	12.559	957	960	962	rBV	804086	915268	16.95%	3.668%
32	12.708	969	973	977	rBV	1193592	1226083	22.71%	4.914%
33	12.776	977	979	982	rVV3	71497	142409	2.64%	0.571%
34	12.891	985	989	991	rVV	177635	267680	4.96%	1.073%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051416\
Data File : BF087068.D
Acq On : 15 May 2016 14:30
Operator : UM/SJ
Sample : H2968-07 2X
Misc :
ALS Vial : 87 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DUPB

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-BF050916.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	12.959	991	995	996	rVV	118840	189536	3.51%	0.760%
36	12.994	996	998	1002	rVV	96448	172617	3.20%	0.692%
37	13.508	1041	1043	1044	rBV	68656	83315	1.54%	0.334%
38	13.748	1061	1064	1069	rBV2	785626	1641452	30.41%	6.579%
39	14.754	1150	1152	1155	rBV	365053	613608	11.37%	2.459%
40	15.062	1177	1179	1182	rBV	263549	495608	9.18%	1.986%
41	15.119	1182	1184	1191	rVB2	446030	749587	13.89%	3.004%

Sum of corrected areas: 24951019

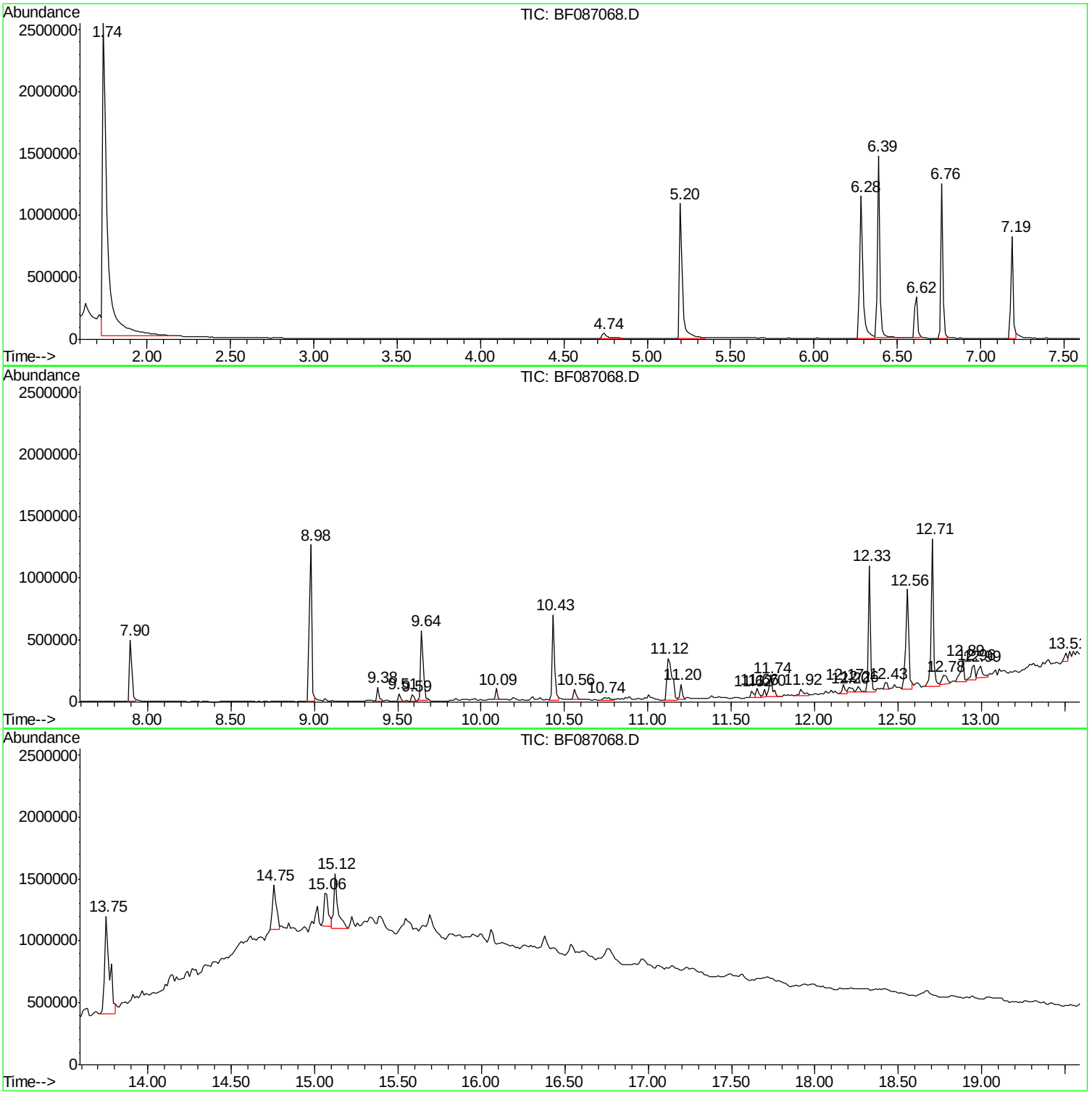
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051416\
Data File : BF087068.D
Acq On : 15 May 2016 14:30
Operator : UM/SJ
Sample : H2968-07 2X
Misc :
ALS Vial : 87 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
DUPB

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.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\BF051416\
Data File : BF087068.D
Acq On : 15 May 2016 14:30
Operator : UM/SJ
Sample : H2968-07 2X
Misc :
ALS Vial : 87 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DUPB

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.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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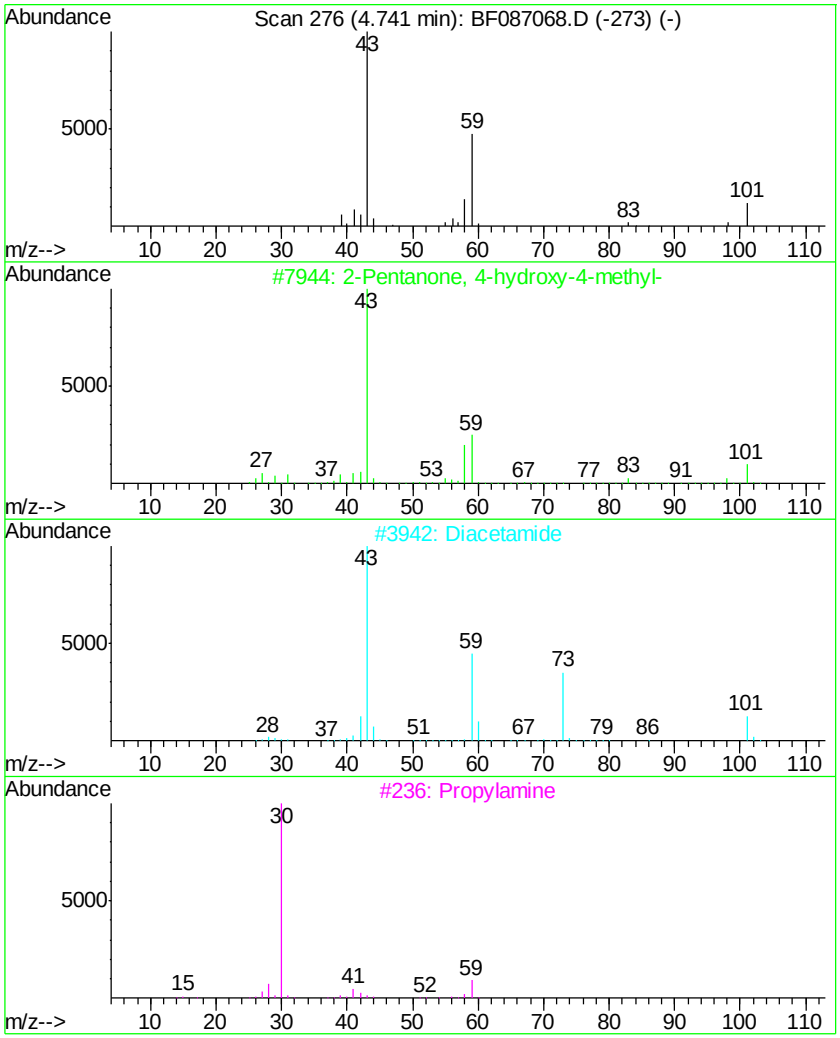
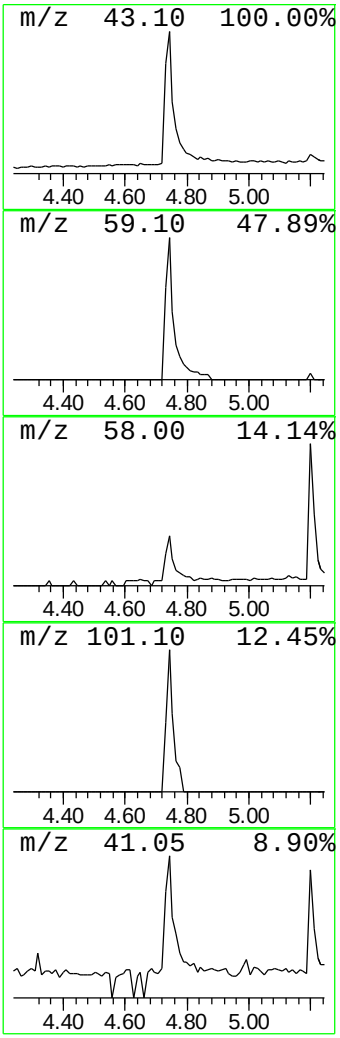
Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.74	4.15 ng	91290	1,4-Dichlorobenzene-d4	6.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2	Diacetamide	101	C4H7NO2	000625-77-4	9
3	Propylamine	59	C3H9N	000107-10-8	9
4	2-Pentanone, 5-(acetyloxy)-	144	C7H12O3	005185-97-7	9
5	5-Aminovaleric acid	117	C5H11NO2	000660-88-8	9





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051416\
Data File : BF087068.D
Acq On : 15 May 2016 14:30
Operator : UM/SJ
Sample : H2968-07 2X
Misc :
ALS Vial : 87 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DUPB

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

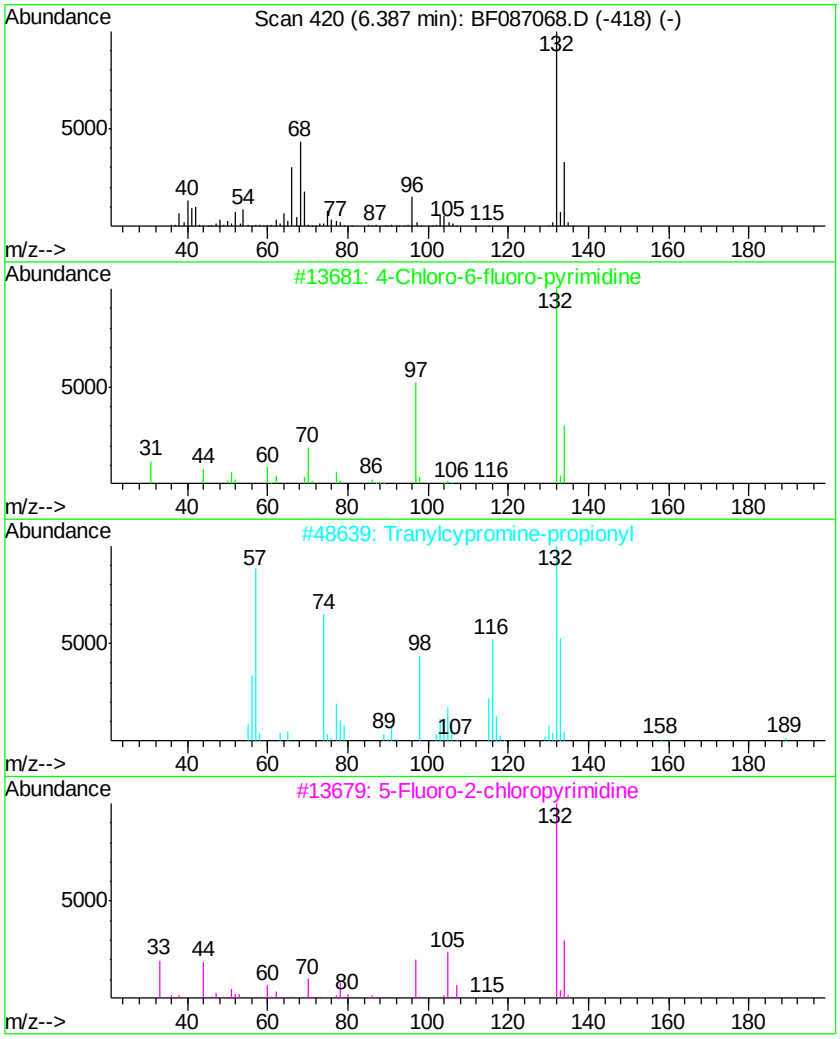
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

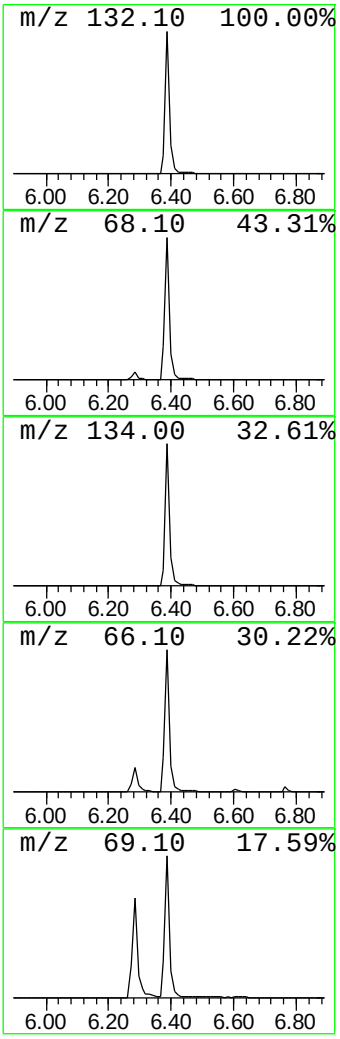
****

Peak Number 3 unknown6.39 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.39	68.67 ng	1511680	1,4-Dichlorobenzene-d4	6.62

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
2	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	22
3	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4	(5-METHYL-2-PYRIDYL)ACETONITRILE	132	C8H8N2	1000241-93-9	10
5	Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051416\
Data File : BF087068.D
Acq On : 15 May 2016 14:30
Operator : UM/SJ
Sample : H2968-07 2X
Misc :
ALS Vial : 87 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DUPB

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.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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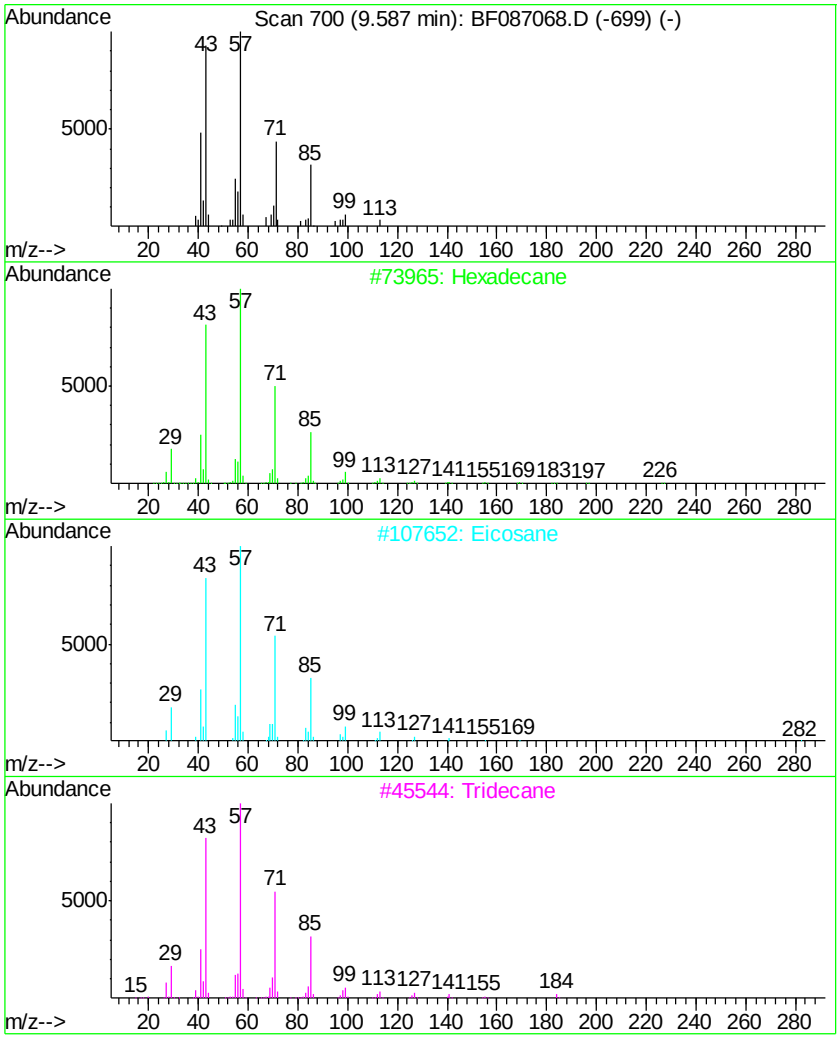
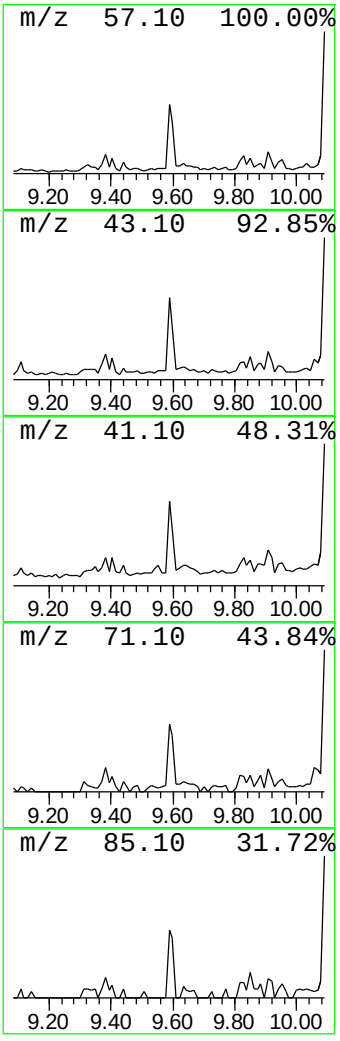
Peak Number 5 Hexadecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.59	2.41 ng	65408	Acenaphthene-d10	9.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	90
2	Eicosane	282	C20H42	000112-95-8	83
3	Tridecane	184	C13H28	000629-50-5	78
4	Pentadecane	212	C15H32	000629-62-9	78
5	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	78





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051416\
Data File : BF087068.D
Acq On : 15 May 2016 14:30
Operator : UM/SJ
Sample : H2968-07 2X
Misc :
ALS Vial : 87 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DUPB

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

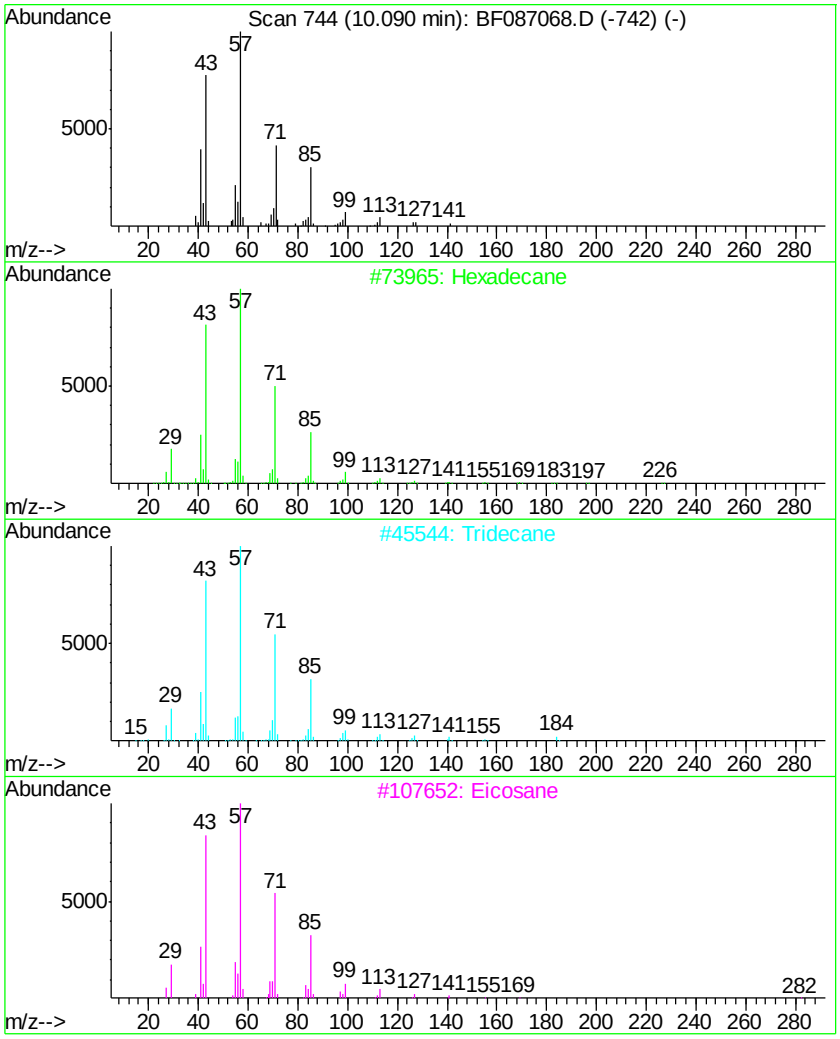
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

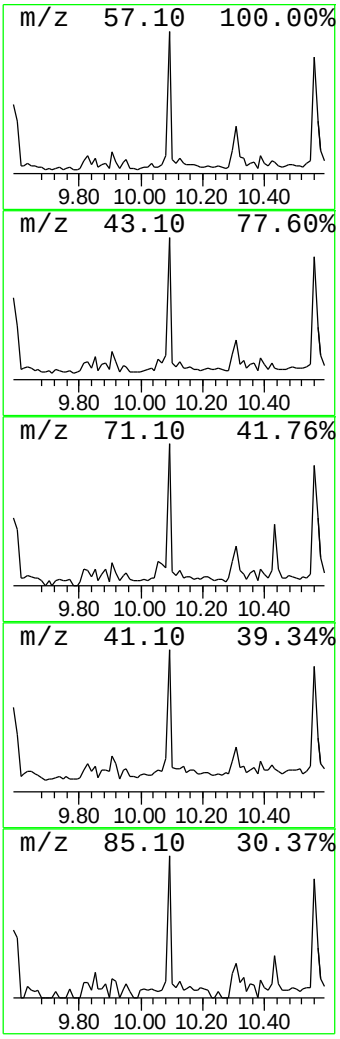
****

Peak Number 6 Tridecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.09	2.74 ng	74317	Acenaphthene-d10	9.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	91
2	Tridecane	184	C13H28	000629-50-5	90
3	Eicosane	282	C20H42	000112-95-8	90
4	Tetradecane	198	C14H30	000629-59-4	72
5	Silane, trichlorooctadecyl-	386	C18H37Cl3Si	000112-04-9	72





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051416\
Data File : BF087068.D
Acq On : 15 May 2016 14:30
Operator : UM/SJ
Sample : H2968-07 2X
Misc :
ALS Vial : 87 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DUPB

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.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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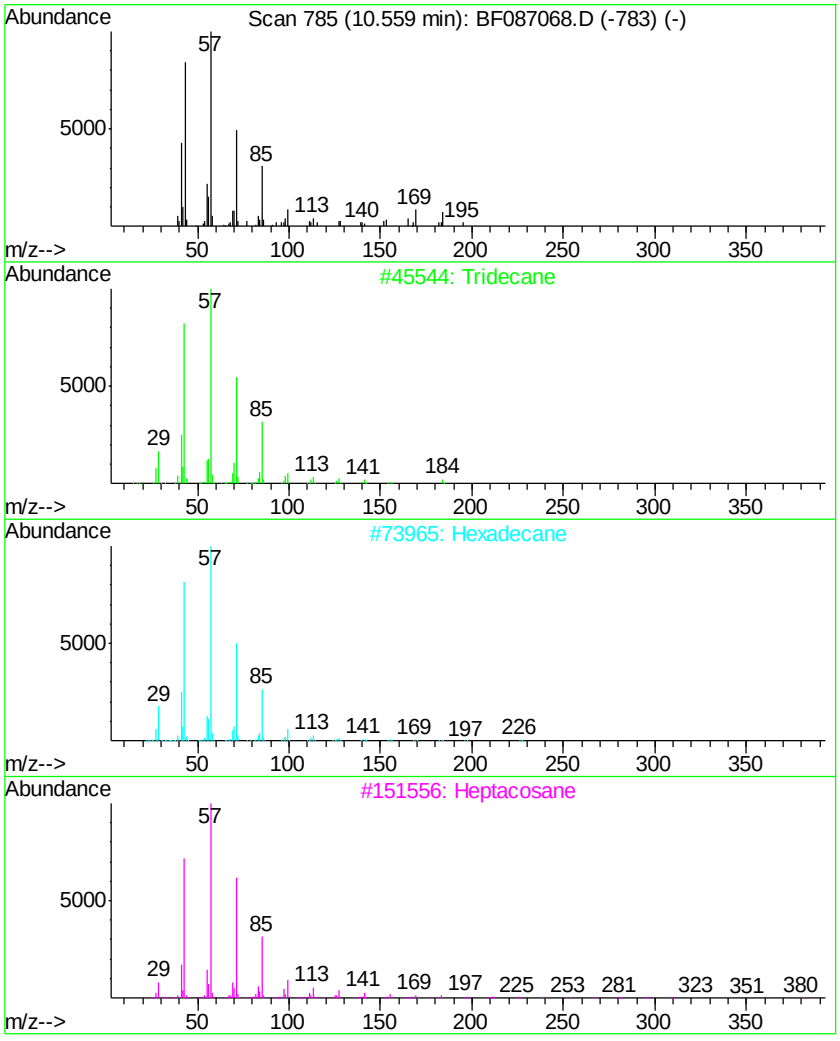
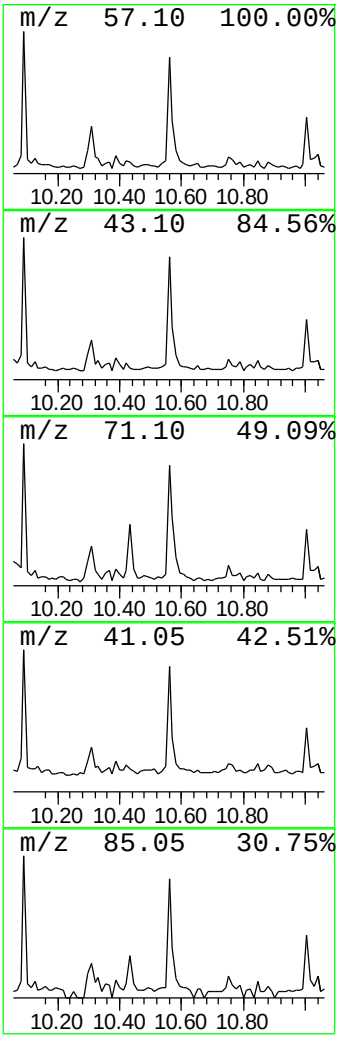
Peak Number 7 Heptacosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.56	2.93 ng	101599	Phenanthrene-d10	11.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	87
2	Hexadecane	226	C16H34	000544-76-3	83
3	Heptacosane	380	C27H56	000593-49-7	80
4	Heptadecane	240	C17H36	000629-78-7	72
5	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	72





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051416\
Data File : BF087068.D
Acq On : 15 May 2016 14:30
Operator : UM/SJ
Sample : H2968-07 2X
Misc :
ALS Vial : 87 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DUPB

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.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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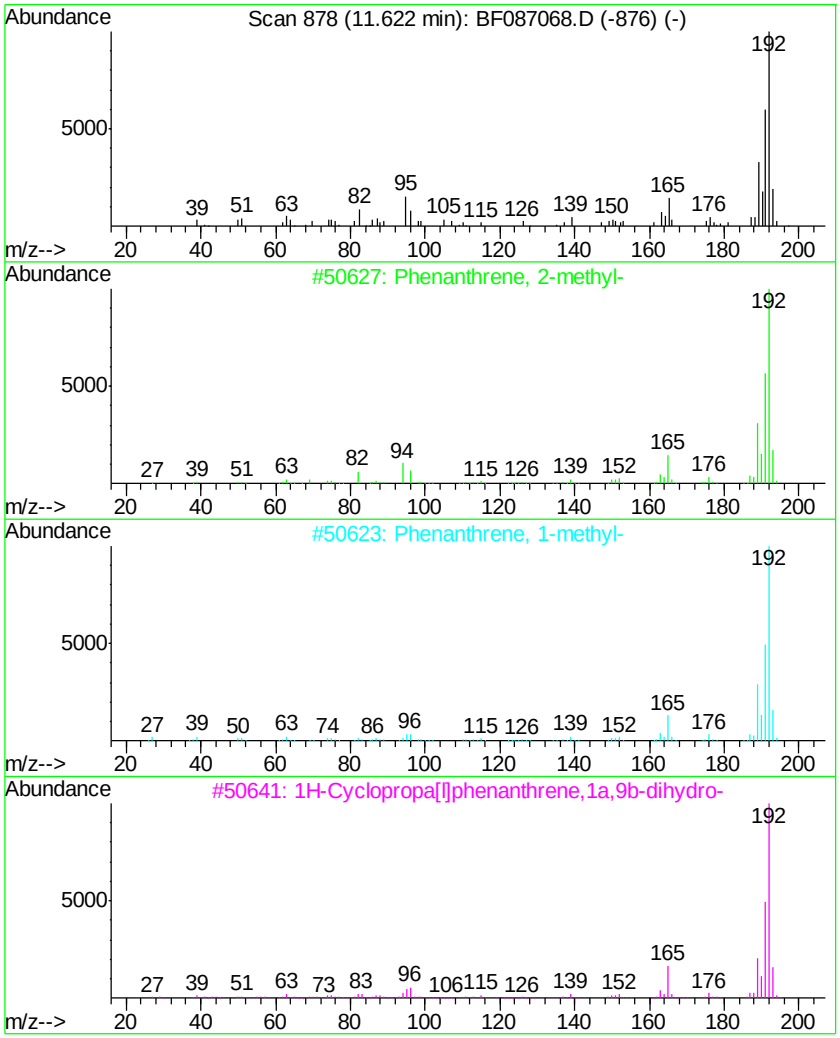
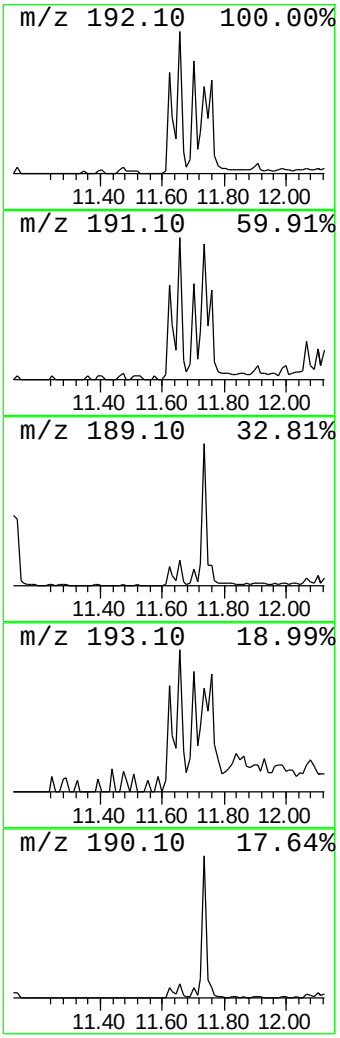
Peak Number 8 Phenanthrene, 2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.62	2.37 ng	82154	Phenanthrene-d10	11.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
3	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	93
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	91
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	91





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051416\
Data File : BF087068.D
Acq On : 15 May 2016 14:30
Operator : UM/SJ
Sample : H2968-07 2X
Misc :
ALS Vial : 87 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DUPB

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.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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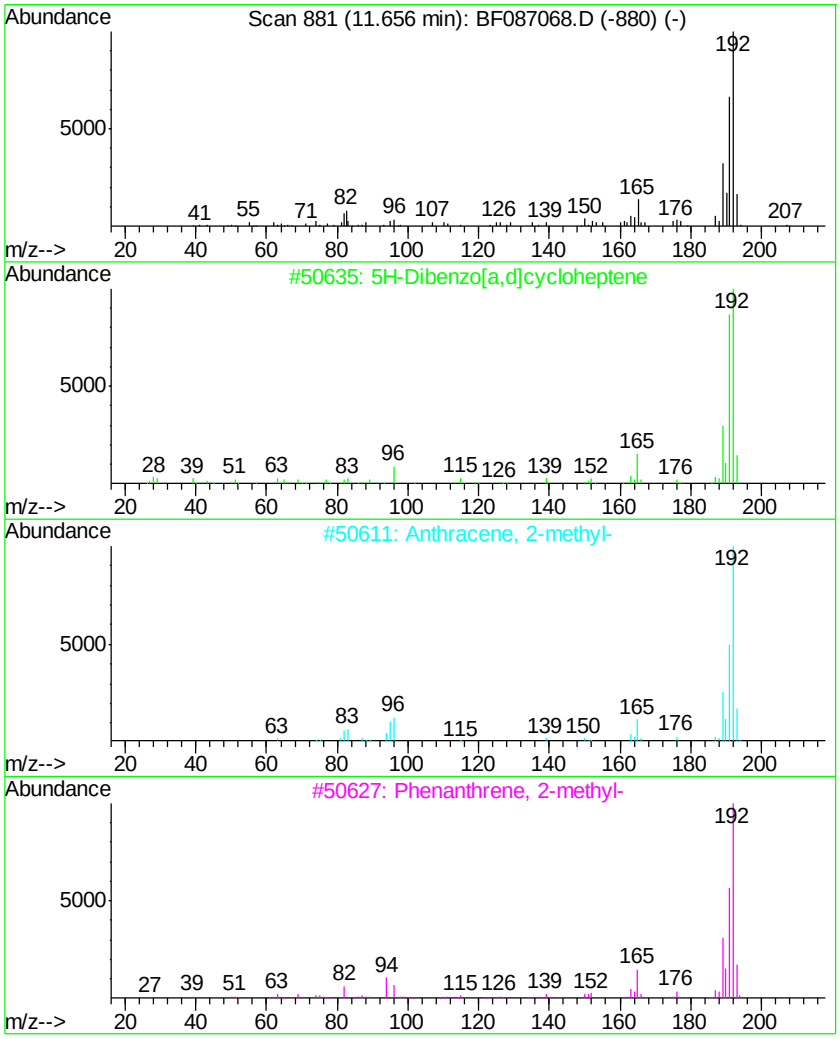
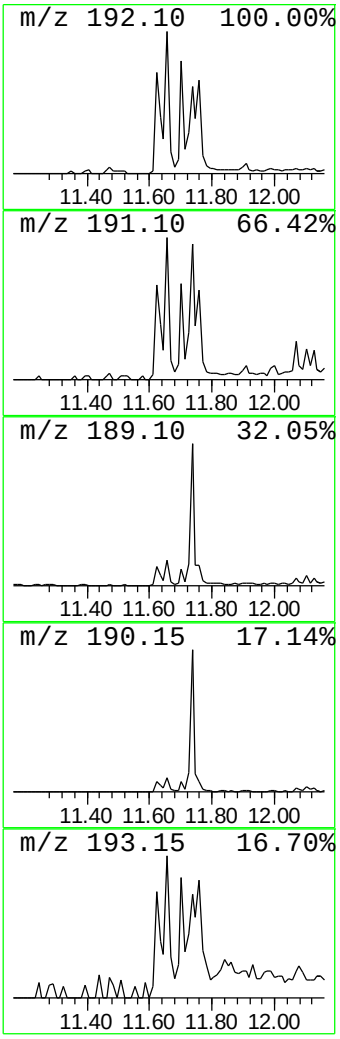
Peak Number 9 5H-Dibenzo[a,d]cycloheptene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.66	2.22 ng	77069	Phenanthrene-d10	11.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	93
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051416\
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Acq On : 15 May 2016 14:30
Operator : UM/SJ
Sample : H2968-07 2X
Misc :
ALS Vial : 87 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DUPB

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.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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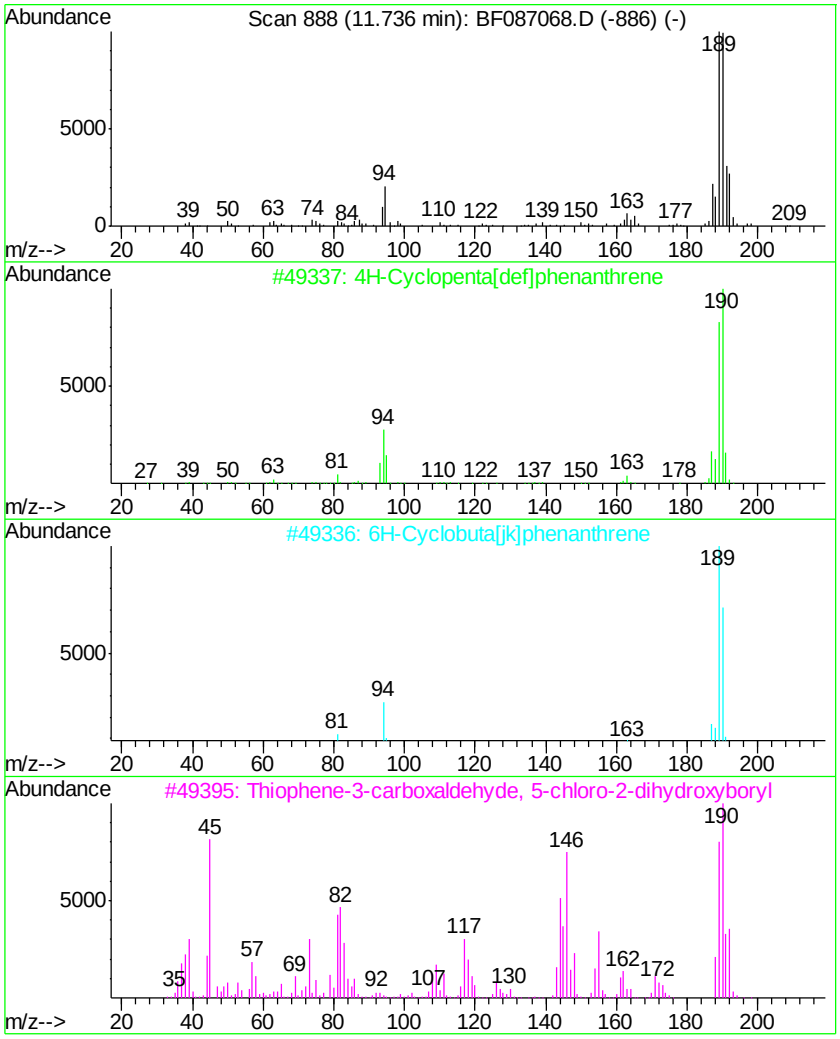
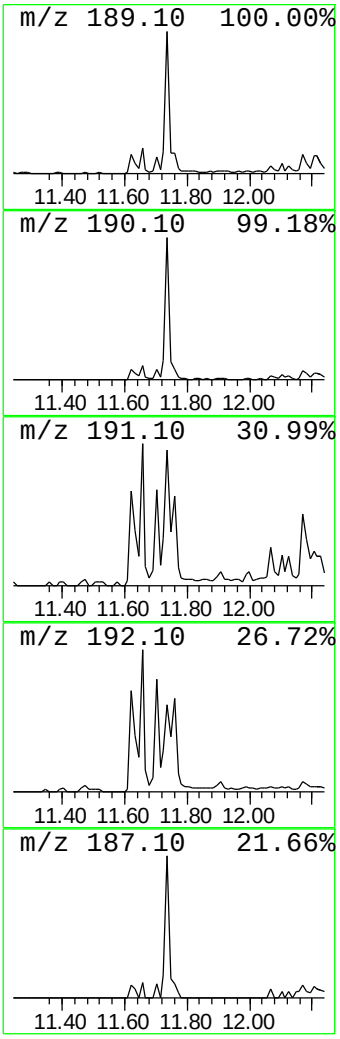
Peak Number 10 4H-Cyclopenta[def]phenanthrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.74	5.95 ng	206527	Phenanthrene-d10	11.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	87
2	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	72
3	Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	52
4	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	45
5	2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	35





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051416\
Data File : BF087068.D
Acq On : 15 May 2016 14:30
Operator : UM/SJ
Sample : H2968-07 2X
Misc :
ALS Vial : 87 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DUPB

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

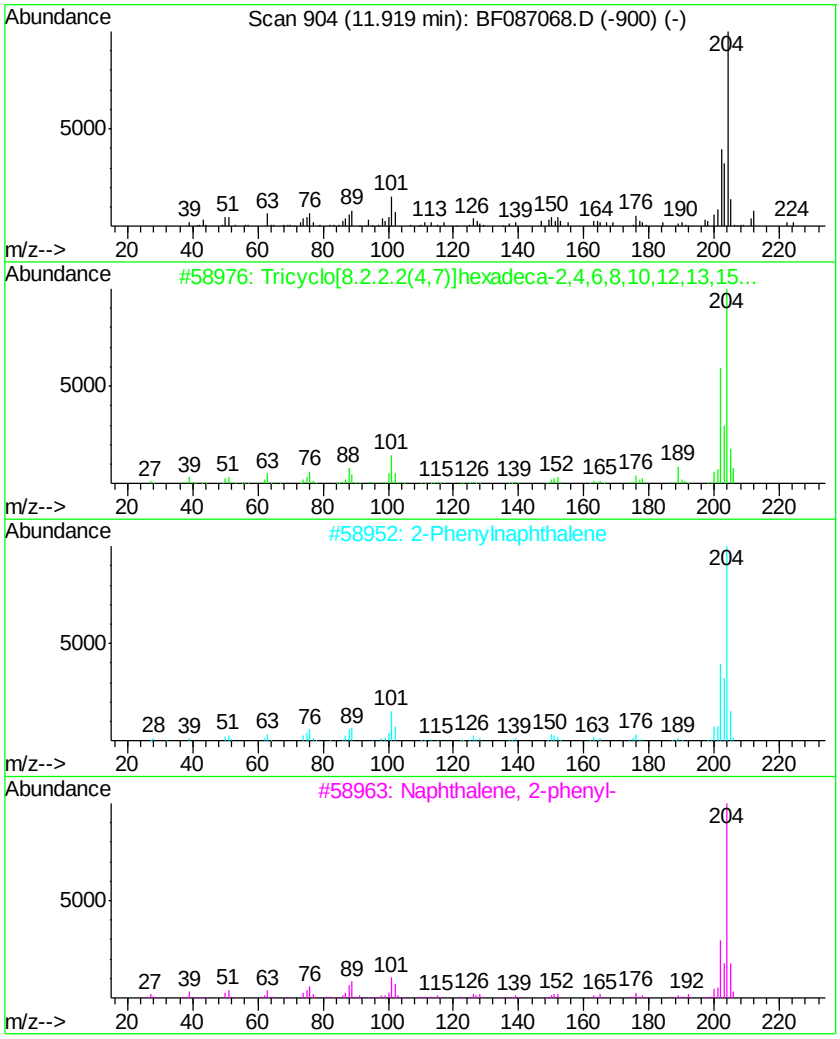
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

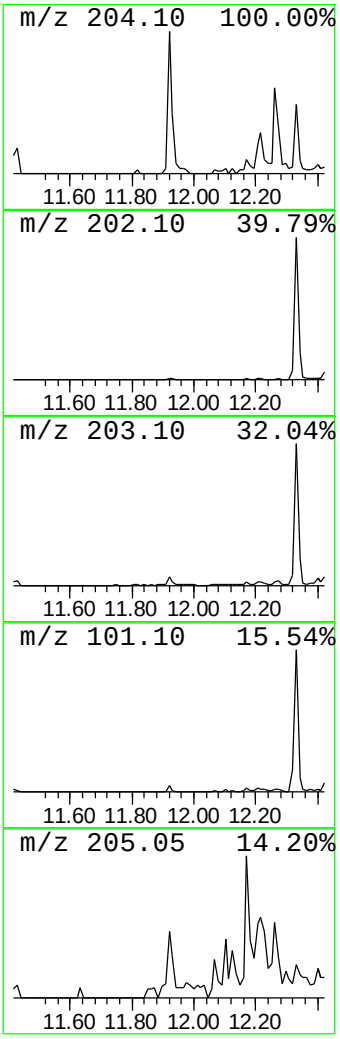
****

Peak Number 11 Tricyclo[8.2.2.2(4,7)]hexad... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.92	2.85 ng	98891	Phenanthrene-d10	11.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	92
2	2-Phenylnaphthalene	204	C16H12	035465-71-5	86
3	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	78
4	1,2,4-Triazolo[2,3-c]thieno[3,2-...	204	C9H8N4S	113246-88-1	49
5	Naphthalene, 1-phenyl-	204	C16H12	000605-02-7	49





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Data File : BF087068.D
Acq On : 15 May 2016 14:30
Operator : UM/SJ
Sample : H2968-07 2X
Misc :
ALS Vial : 87 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DUPB

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

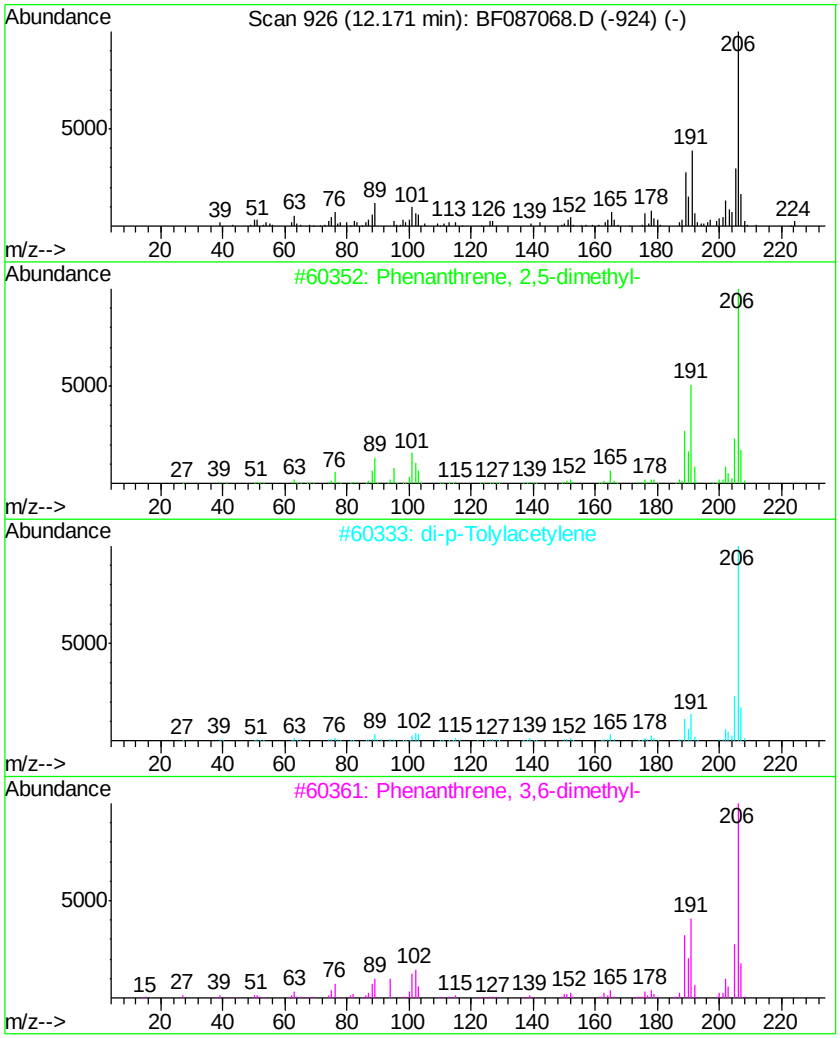
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

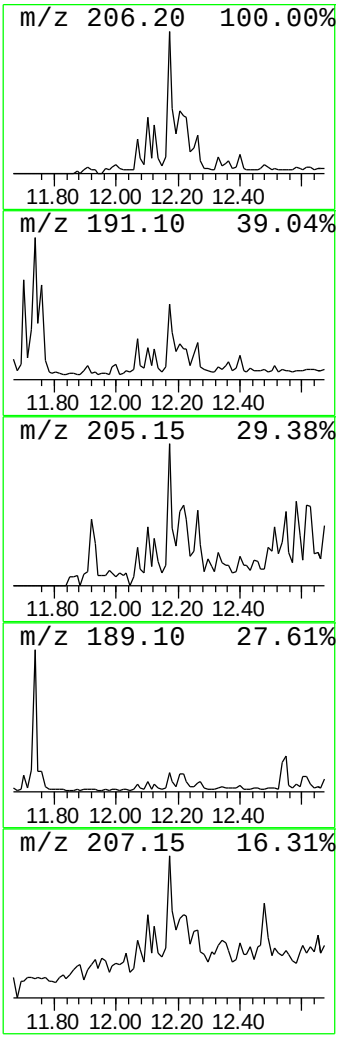
****

Peak Number 12 Phenanthrene, 2,5-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.17	2.57 ng	89121	Phenanthrene-d10	11.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	94
2	di-p-Tolylacetylene	206	C16H14	002789-88-0	94
3	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	93
4	Naphthalene, 1,2-dihydro-4-phenyl-	206	C16H14	007469-40-1	93
5	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	90





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Acq On : 15 May 2016 14:30
Operator : UM/SJ
Sample : H2968-07 2X
Misc :
ALS Vial : 87 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DUPB

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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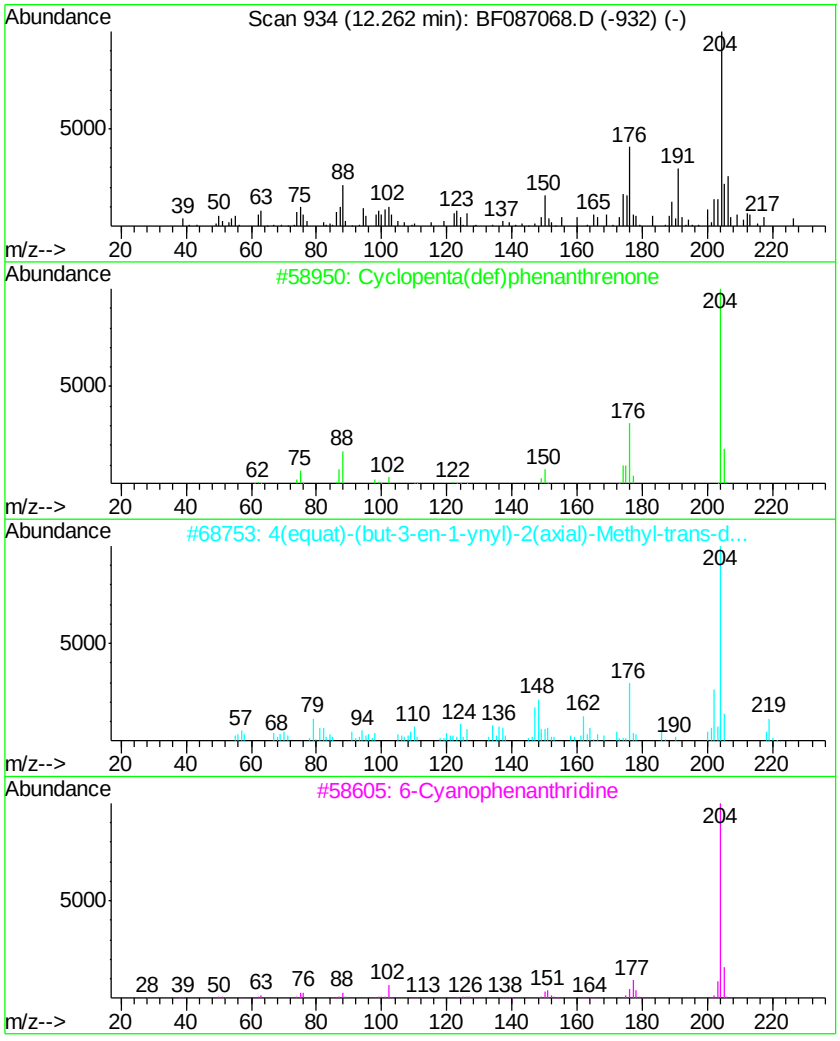
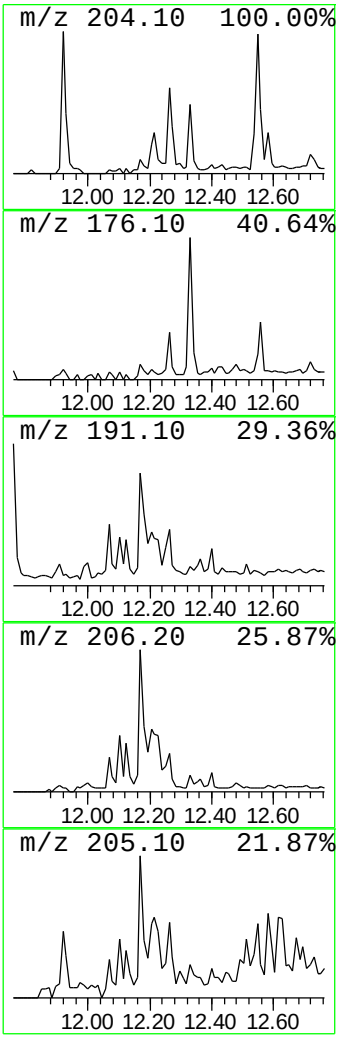
Peak Number 13 Cyclopenta(def)phenanthrenone Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.26	2.06 ng	71573	Phenanthrene-d10	11.13

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	93
2	4(equat)-(but-3-en-1-ynyl)-2(axi...	219	C14H21NO	038423-22-2	47
3	6-Cyanophenanthridine	204	C14H8N2	002176-28-5	43
4	1-Methyl-4-ethyl 2-phenylsuccinate	236	C13H16O4	132545-36-9	43
5	9-Acridinecarbonitrile	204	C14H8N2	005326-19-2	43





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Acq On : 15 May 2016 14:30
Operator : UM/SJ
Sample : H2968-07 2X
Misc :
ALS Vial : 87 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DUPB

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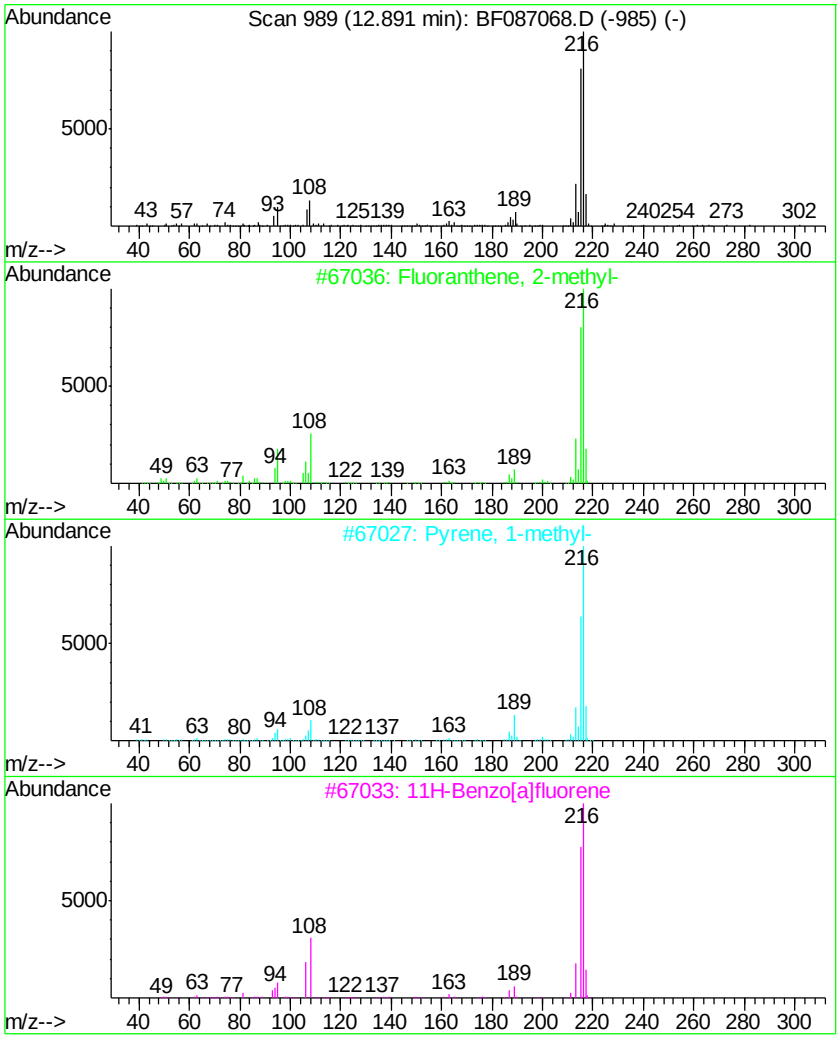
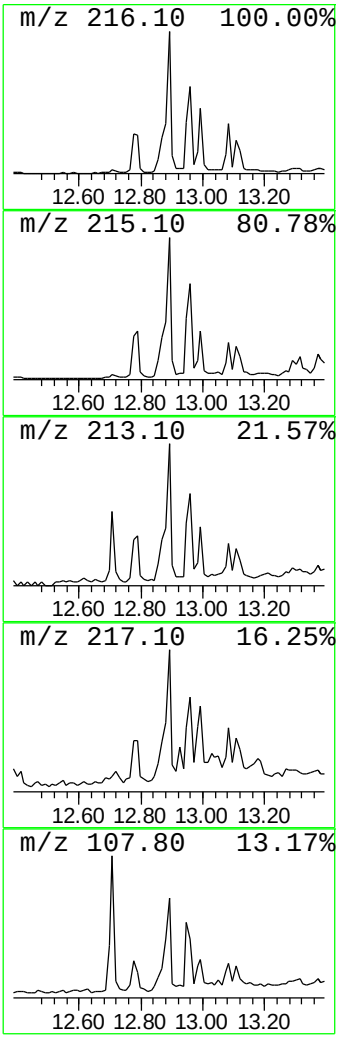
Peak Number 15 Fluoranthene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.89	3.26 ng	267680	Chrysene-d12	13.76

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	93
2	Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	93
4	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	91
5	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	83





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Acq On : 15 May 2016 14:30
Operator : UM/SJ
Sample : H2968-07 2X
Misc :
ALS Vial : 87 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DUPB

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.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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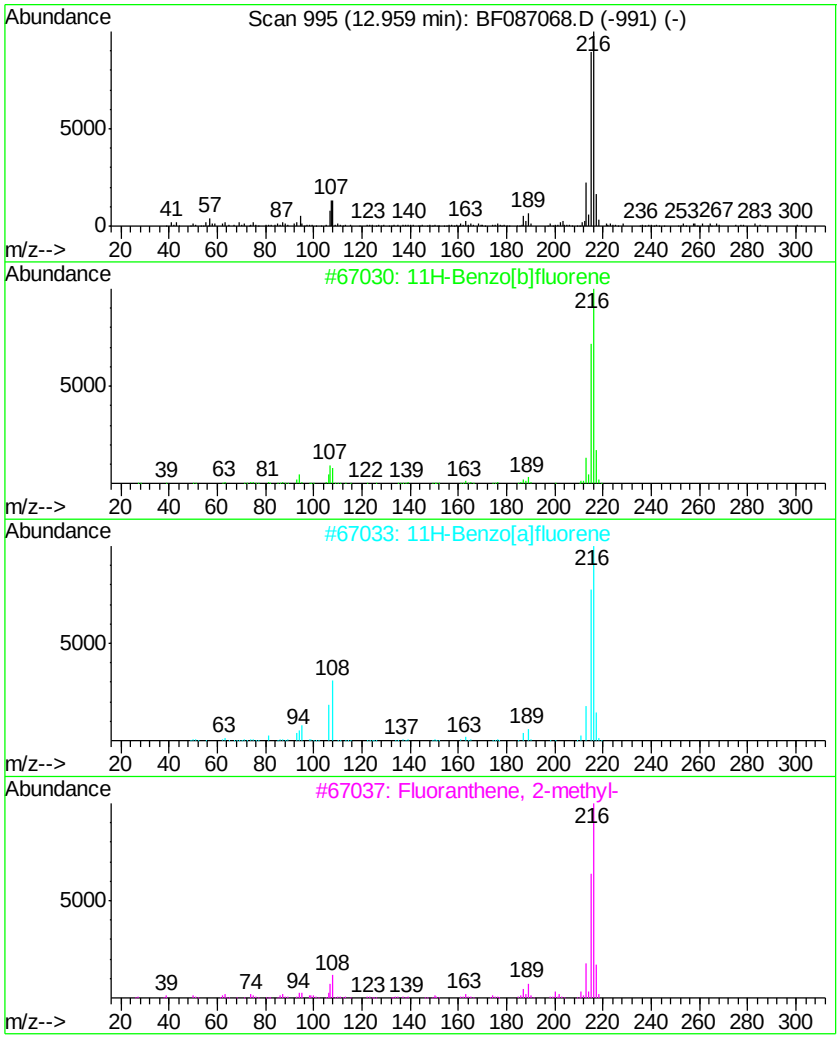
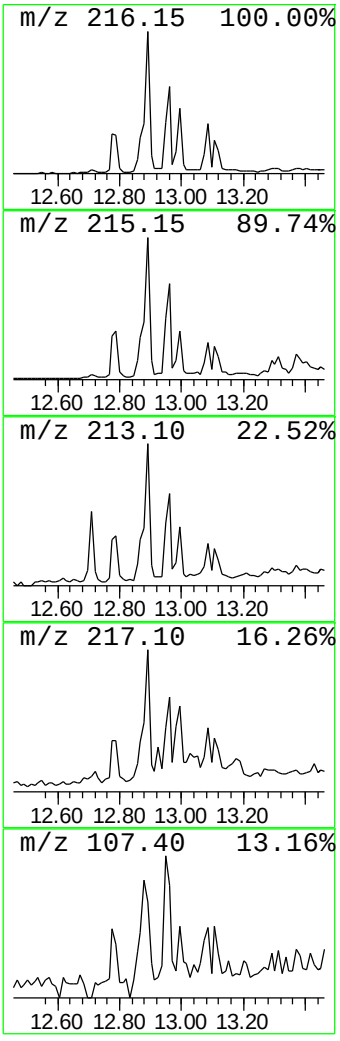
Peak Number 16 11H-Benzo[b]fluorene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.96	2.31 ng	189536	Chrysene-d12	13.76

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	96
2	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
3	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	80
4	Pyrene, 1-methyl-	216	C17H12	002381-21-7	68
5	1,3,5-Triazine-2,4,6-triamine, N...	216	C10H12N6	046731-79-7	59





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Acq On : 15 May 2016 14:30
Operator : UM/SJ
Sample : H2968-07 2X
Misc :
ALS Vial : 87 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DUPB

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.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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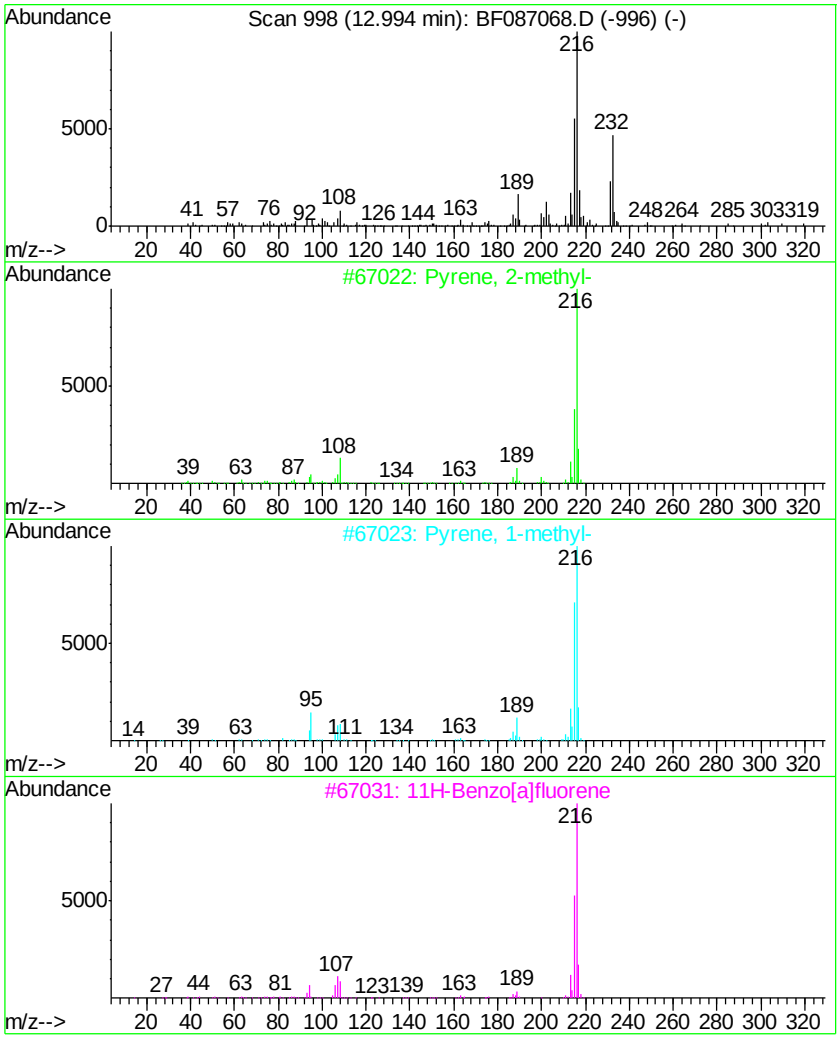
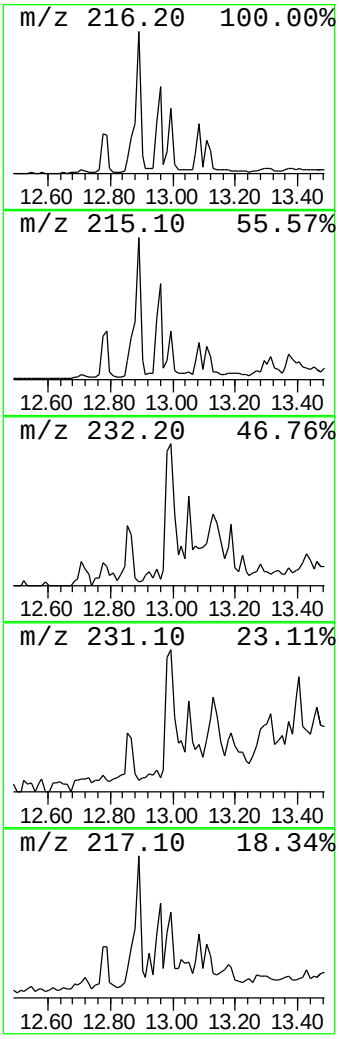
Peak Number 17 Pyrene, 2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.99	2.10 ng	172617	Chrysene-d12	13.76

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 2-methyl-	216	C17H12	003442-78-2	90
2	Pyrene, 1-methyl-	216	C17H12	002381-21-7	64
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	64
4	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	60
5	Pyrene, 4-methyl-	216	C17H12	003353-12-6	55





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051416\
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Acq On : 15 May 2016 14:30
Operator : UM/SJ
Sample : H2968-07 2X
Misc :
ALS Vial : 87 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : DUPB

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050916.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
2-Pentanone, 4-hy...	4.74	4.2	ng	91290	1	6.62	440253	20.0
unknown6.39	6.39	68.7	ng	1511680	1	6.62	440253	20.0
Hexadecane	9.59	2.4	ng	65408	3	9.64	543151	20.0
Tridecane	10.09	2.7	ng	74317	3	9.64	543151	20.0
Heptacosane	10.56	2.9	ng	101599	4	11.13	694073	20.0
Phenanthrene, 2-m...	11.62	2.4	ng	82154	4	11.13	694073	20.0
5H-Dibenzo[a,d]cy...	11.66	2.2	ng	77069	4	11.13	694073	20.0
4H-Cyclopenta[def...	11.74	6.0	ng	206527	4	11.13	694073	20.0
Tricyclo[8.2.2.2(...	11.92	2.9	ng	98891	4	11.13	694073	20.0
Phenanthrene, 2,5...	12.17	2.6	ng	89121	4	11.13	694073	20.0
Cyclopenta(def)ph...	12.26	2.1	ng	71573	4	11.13	694073	20.0
Fluoranthene, 2-m...	12.89	3.3	ng	267680	5	13.76	1641450	20.0
11H-Benzo[b]fluorene	12.96	2.3	ng	189536	5	13.76	1641450	20.0
Pyrene, 2-methyl-	12.99	2.1	ng	172617	5	13.76	1641450	20.0