LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011317\
Data File : BF092271.D
Acq On : 14 Jan 2017 4:58
Operator : UM/SJ
Sample : I1182-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TOD-BC-0.2

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	1.890	7	9	28	rVB	44099	234689	7.06%	0.610%
2	3.970	188	191	197	rBV	20414	41848	1.26%	0.109%
3	4.701	252	255	264	rBV	1125041	1594539	47.95%	4.145%
4	5.170	294	296	305	rBV	2925406	3325582	100.00%	8.644%
5	6.210	384	387	388	rBV	267091	280027	8.42%	0.728%
6	6.244	388	390	397	rVV	3046690	3200221	96.23%	8.319%
7	6.347	397	399	402	rVB	3666235	3260906	98.06%	8.476%
8	6.576	416	419	421	rBV	1160020	1114332	33.51%	2.897%
9	6.724	430	432	435	rBV	1817500	2292023	68.92%	5.958%
10	7.147	466	469	471	rBV	2189099	2116868	63.65%	5.503%
11	7.856	529	531	534	rBV	927976	1305102	39.24%	3.392%
12	8.439	580	582	584	rVB	64563	68969	2.07%	0.179%
13	8.942	623	626	628	rBV	3114037	2725514	81.96%	7.085%
14	9.079	636	638	639	rVB	34716	39259	1.18%	0.102%
15	9.342	659	661	664	rBV	501107	417866	12.57%	1.086%
16	9.605	682	684	686	rBV	849029	1082258	32.54%	2.813%
17	9.639	686	687	691	rVB2	55194	68562	2.06%	0.178%
18	10.153	725	732	735	rVB2	40571	52737	1.59%	0.137%
19	10.393	751	753	756	rBV2	1065977	1438559	43.26%	3.739%
20	10.530	763	765	769	rVB	63512	69255	2.08%	0.180%
21	10.976	802	804	808	rVB2	72899	98017	2.95%	0.255%
22	11.091	811	814	818	rVV2	770212	1283196	38.59%	3.336%
23	11.159	818	820	822	rVV	114613	116564	3.51%	0.303%
24	11.331	833	835	839	rBV	279606	360698	10.85%	0.938%
25	11.399	839	841	848	rVB	71532	101138	3.04%	0.263%
26	11.582	852	857	859	rBV	56377	130973	3.94%	0.340%
27	11.616	859	860	861	rVV	75730	56262	1.69%	0.146%
28	11.639	861	862	865	rVV2	141985	180456	5.43%	0.469%
29	11.696	865	867	872	rVB2	116307	172326	5.18%	0.448%
30	11.811	875	877	879	rVB	49336	77757	2.34%	0.202%
31	11.879	881	883	885	rBV3	37574	50028	1.50%	0.130%
32	12.153	904	907	910	rBV2	568796	853122	25.65%	2.218%
33	12.291	917	919	921	rBV	740125	682547	20.52%	1.774%
34	12.359	921	925	926	rVV3	29436	57273	1.72%	0.149%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011317\
Data File : BF092271.D
Acq On : 14 Jan 2017 4:58
Operator : UM/SJ
Sample : I1182-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TOD-BC-0.2

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	12.439	929	932	936	rBV3	44472	106898	3.21%	0.278%
36	12.519	936	939	941	rBV	592503	747076	22.46%	1.942%
37	12.565	941	943	948	rVB3	56120	92513	2.78%	0.240%
38	12.668	950	952	955	rVB	1690875	1823005	54.82%	4.739%
39	12.748	955	959	960	rBV3	34896	51224	1.54%	0.133%
40	12.851	964	968	970	rVB	85793	151709	4.56%	0.394%
41	12.919	970	974	975	rBV	89448	140758	4.23%	0.366%
42	12.954	975	977	981	rVB2	74066	107024	3.22%	0.278%
43	13.045	983	985	986	rBV	42876	49473	1.49%	0.129%
44	13.079	986	988	991	rBV3	30864	58773	1.77%	0.153%
45	13.125	991	992	995	rBV2	54243	85522	2.57%	0.222%
46	13.216	998	1000	1002	rBV2	84984	109782	3.30%	0.285%
47	13.354	1008	1012	1015	rBV2	49484	112164	3.37%	0.292%
48	13.468	1019	1022	1023	rBV	64141	92942	2.79%	0.242%
49	13.582	1029	1032	1035	rBV2	141262	237423	7.14%	0.617%
50	13.708	1040	1043	1048	rBV2	929371	1700485	51.13%	4.420%
51	13.811	1050	1052	1054	rBV2	41928	65133	1.96%	0.169%
52	14.097	1075	1077	1079	rBV	46953	72593	2.18%	0.189%
53	14.211	1084	1087	1091	rVV3	112823	269008	8.09%	0.699%
54	14.565	1116	1118	1120	rBV2	121887	151447	4.55%	0.394%
55	14.714	1128	1131	1135	rBV	244124	521548	15.68%	1.356%
56	14.794	1135	1138	1140	rBV2	382696	493763	14.85%	1.283%
57	14.965	1151	1153	1155	rVB	133249	154229	4.64%	0.401%
58	15.011	1155	1157	1160	rVV	170942	256451	7.71%	0.667%
59	15.068	1160	1162	1168	rVB	610708	832784	25.04%	2.165%
60	15.388	1189	1190	1193	rVB3	80560	76654	2.30%	0.199%
61	15.468	1195	1197	1200	rBV	312217	470671	14.15%	1.223%
62	16.302	1268	1270	1273	rVB	83829	145943	4.39%	0.379%
63	16.668	1300	1302	1307	rBV	57905	109522	3.29%	0.285%
64	16.977	1327	1329	1333	rBV4	75407	234713	7.06%	0.610%

Sum of corrected areas: 38470703

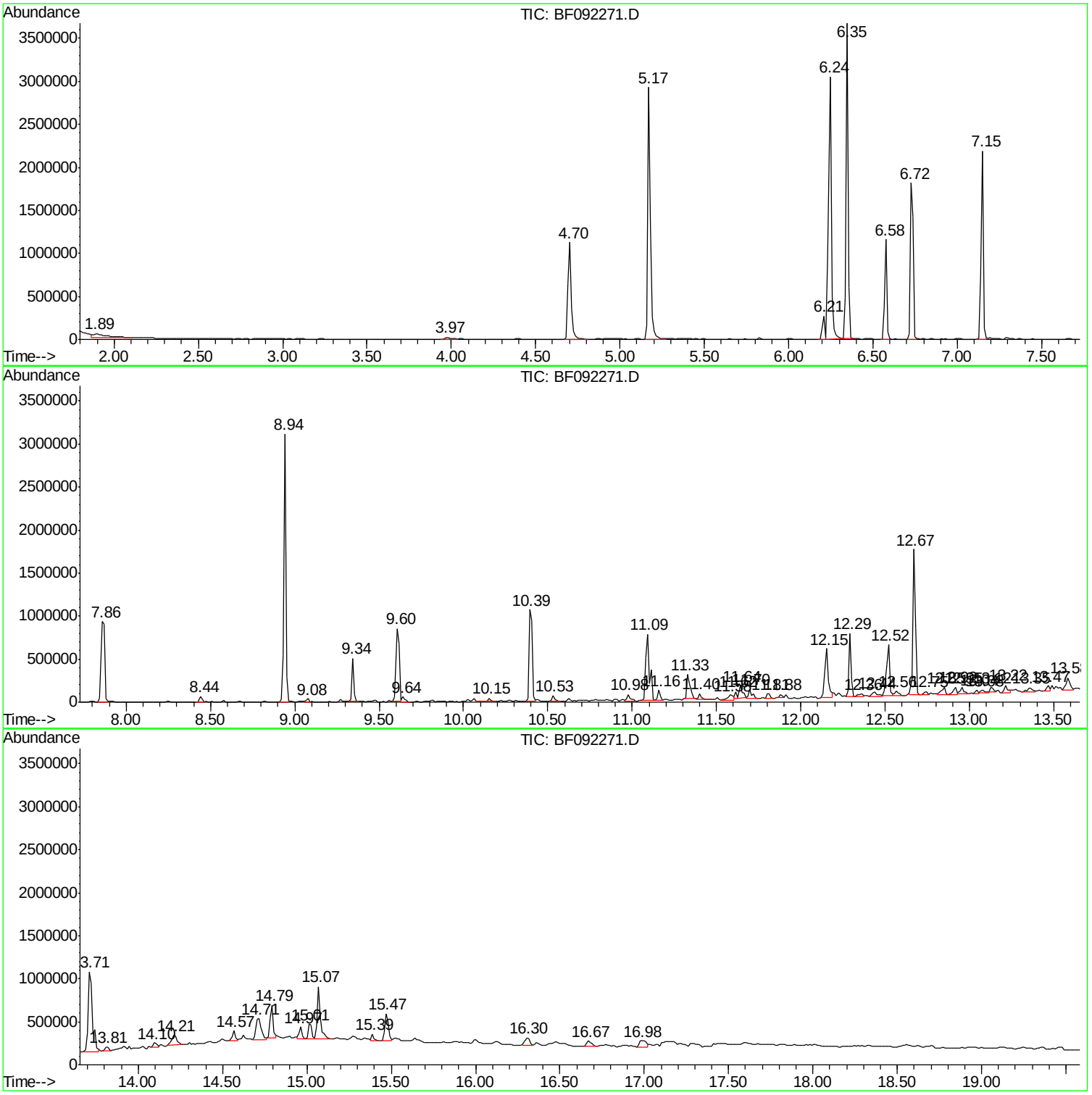
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011317\
Data File : BF092271.D
Acq On : 14 Jan 2017 4:58
Operator : UM/SJ
Sample : I1182-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TOD-BC-0.2

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\BF011317\
Data File : BF092271.D
Acq On : 14 Jan 2017 4:58
Operator : UM/SJ
Sample : I1182-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TOD-BC-0.2

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

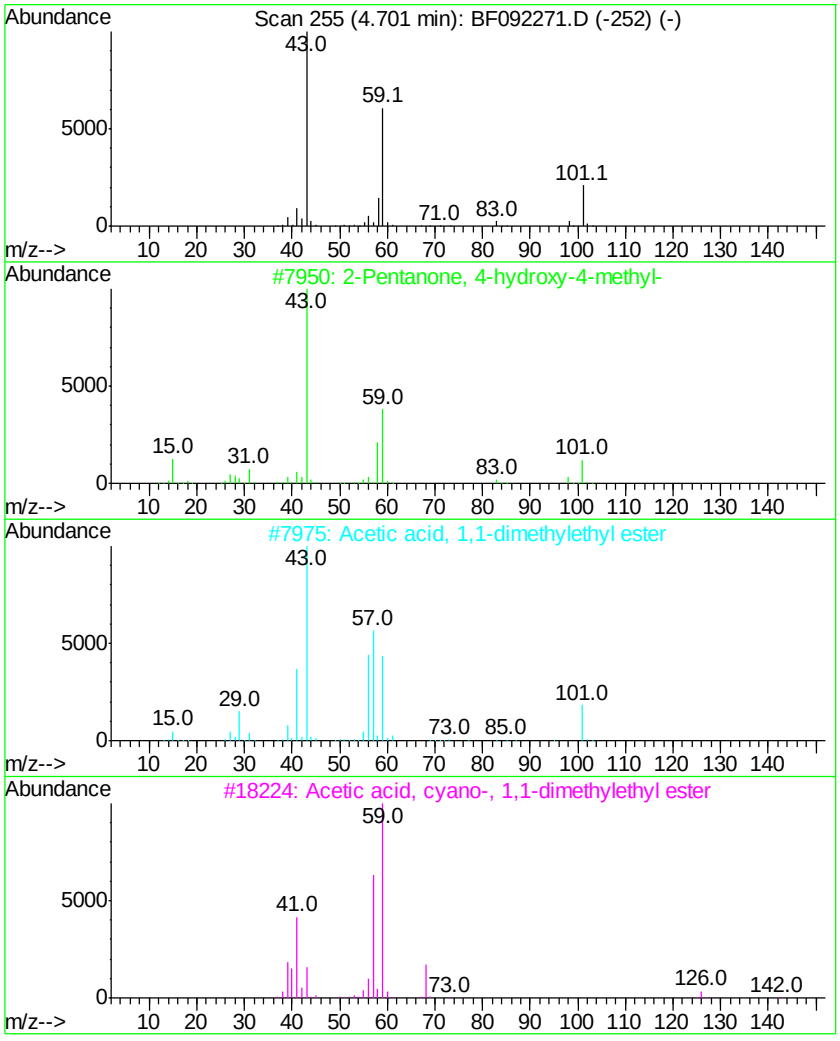
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

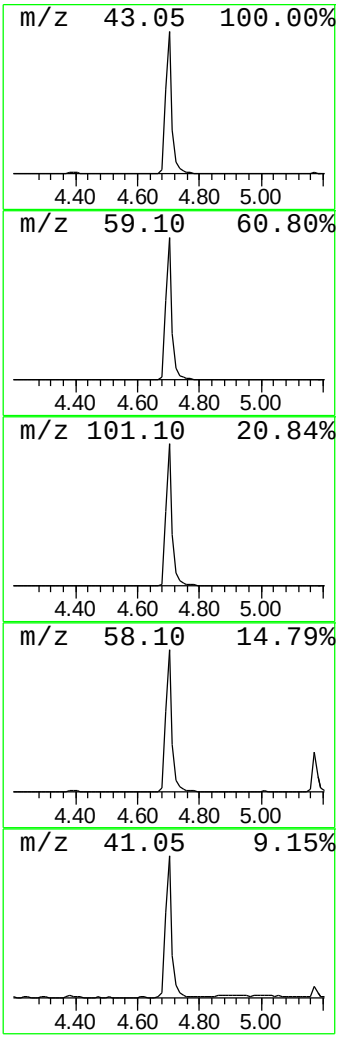
****

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.70	28.62 ng	1594540	1,4-Dichlorobenzene-d4	6.58

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	39
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
4	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF011317\
Data File : BF092271.D
Acq On : 14 Jan 2017 4:58
Operator : UM/SJ
Sample : I1182-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TOD-BC-0.2

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

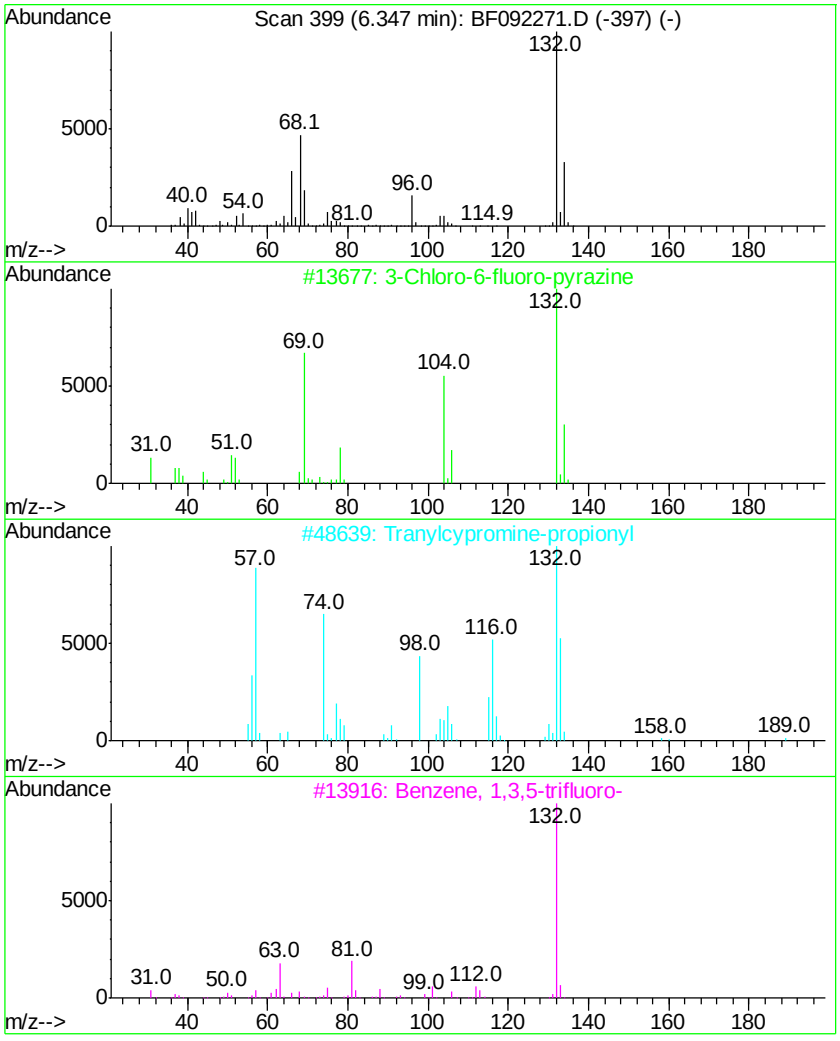
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

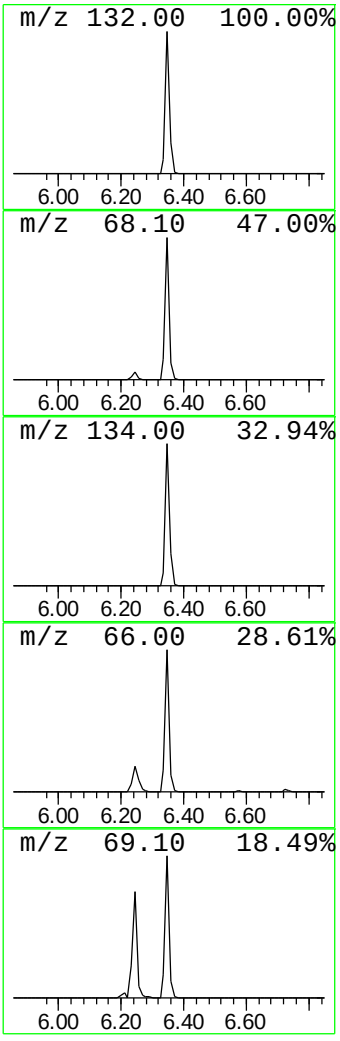
****

Peak Number 3 unknown6.35 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.35	58.53 ng	3260910	1,4-Dichlorobenzene-d4	6.58

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3	Benzene, 1,3,5-trifluoro-	132	C6H3F3	000372-38-3	9
4	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011317\
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Acq On : 14 Jan 2017 4:58
Operator : UM/SJ
Sample : I1182-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TOD-BC-0.2

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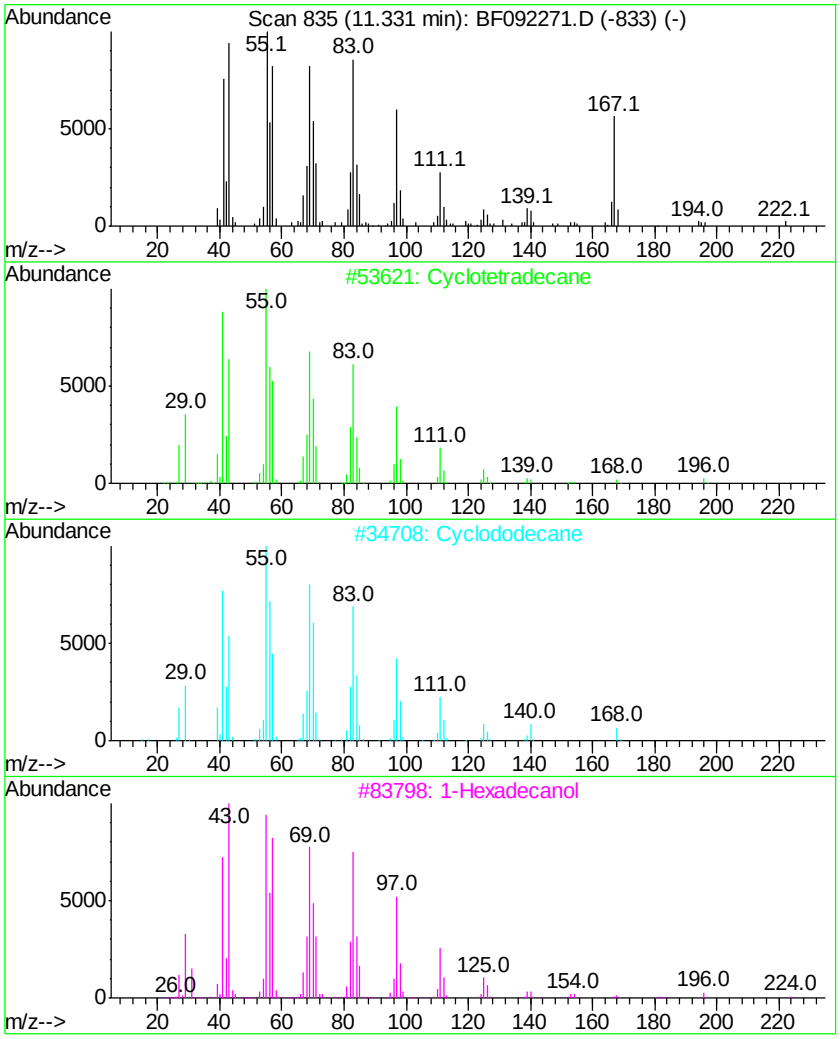
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

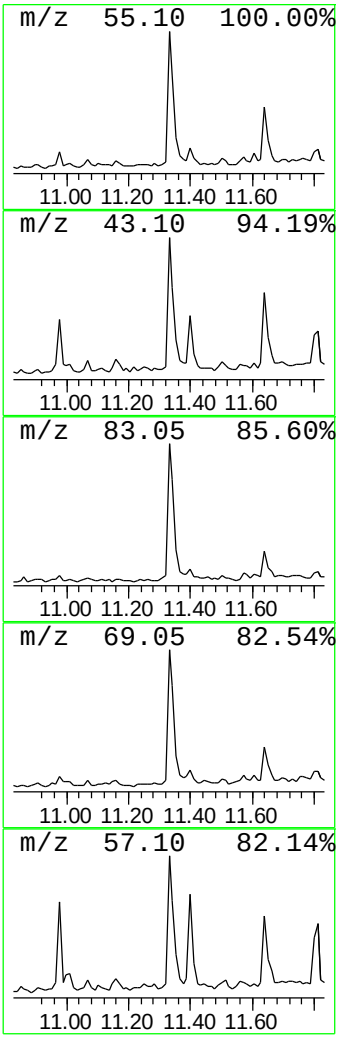
****

Peak Number 5 Cyclotetradecane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.33	5.62 ng	360698	Phenanthrene-d10	11.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclotetradecane	196	C14H28	000295-17-0	96
2	Cyclododecane	168	C12H24	000294-62-2	92
3	1-Hexadecanol	242	C16H34O	036653-82-4	91
4	2-Tetradecene, (E)-	196	C14H28	035953-53-8	89
5	1-Tetradecene	196	C14H28	001120-36-1	87





Data Path : Z:\HPCHEM1\BNA F\DATA\BF011317\
Data File : BF092271.D
Acq On : 14 Jan 2017 4:58
Operator : UM/SJ
Sample : I1182-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TOD-BC-0.2

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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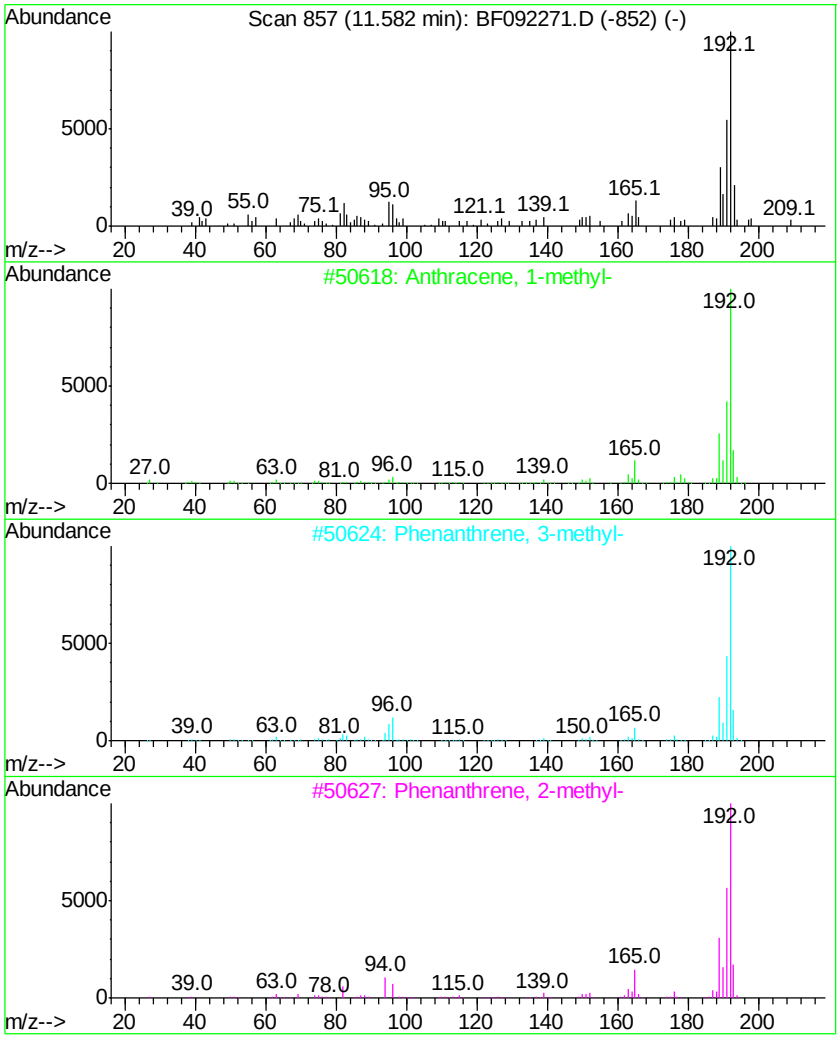
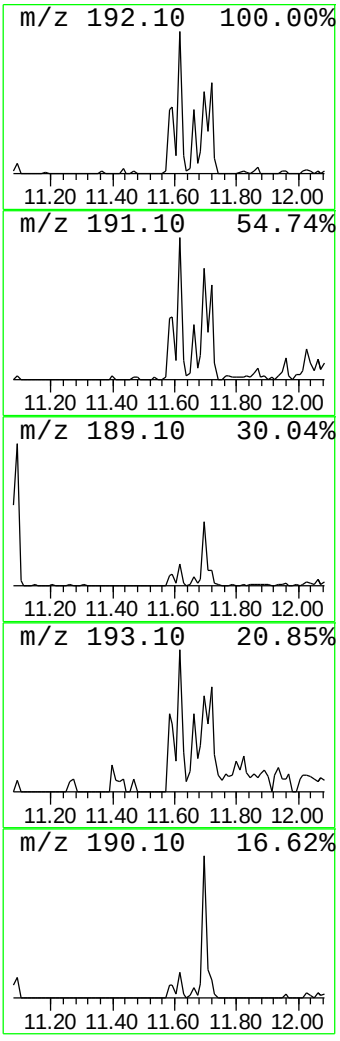
Peak Number 6 Anthracene, 1-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.58	2.04 ng	130973	Phenanthrene-d10	11.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 1-methyl-	192	C15H12	000610-48-0	94
2	Phenanthrene, 3-methyl-	192	C15H12	000832-71-3	93
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	90
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90





Data Path : Z:\HPCHEM1\BNA F\DATA\BF011317\
Data File : BF092271.D
Acq On : 14 Jan 2017 4:58
Operator : UM/SJ
Sample : I1182-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TOD-BC-0.2

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

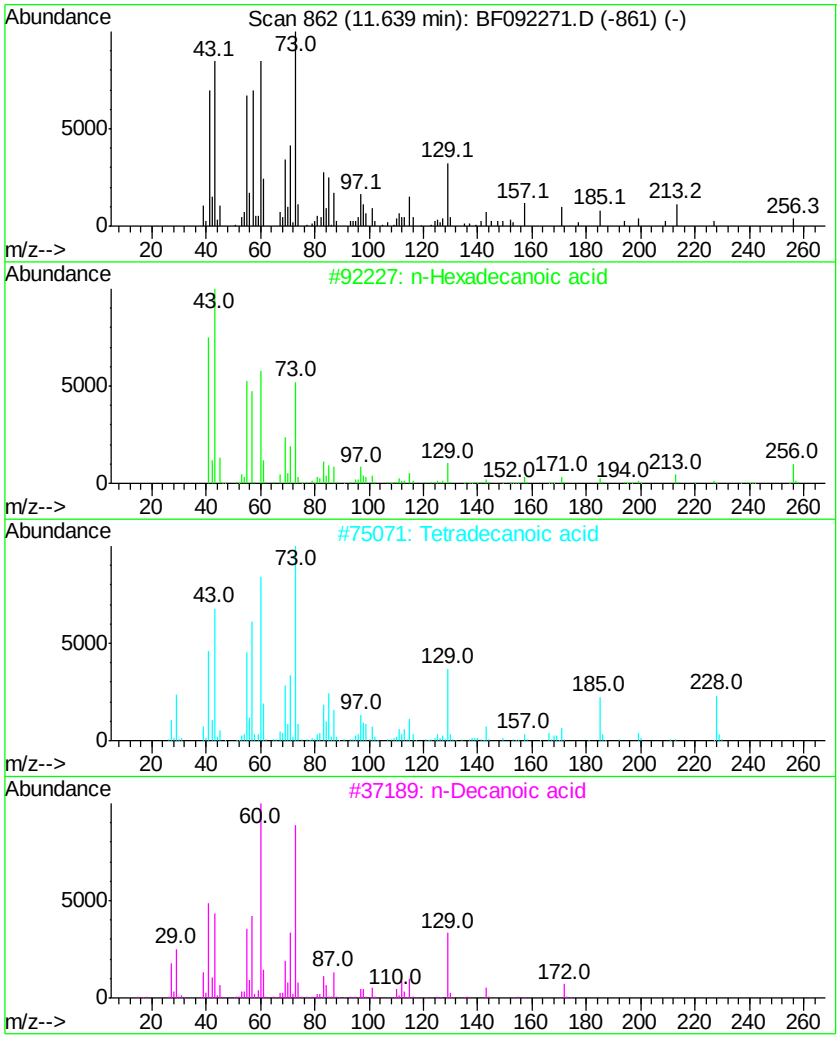
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

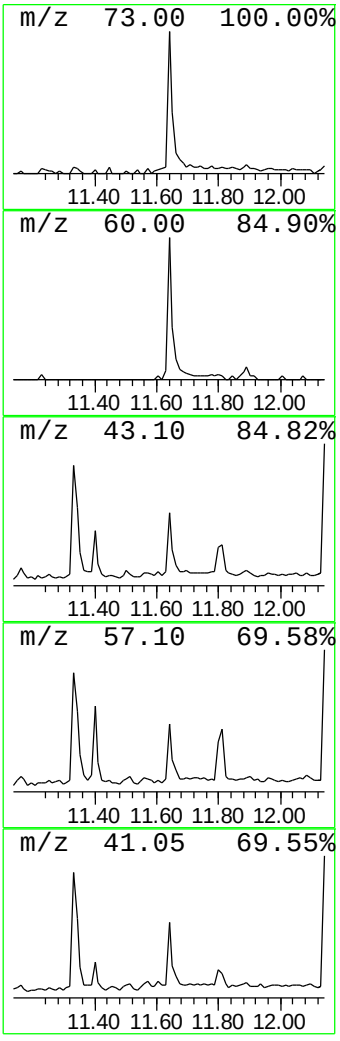
****

Peak Number 7 n-Hexadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.64	2.81 ng	180456	Phenanthrene-d10	11.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	83
3	n-Decanoic acid	172	C10H20O2	000334-48-5	64
4	Tridecanoic acid	214	C13H26O2	000638-53-9	64
5	Dodecanoic acid	200	C12H24O2	000143-07-7	58





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011317\
Data File : BF092271.D
Acq On : 14 Jan 2017 4:58
Operator : UM/SJ
Sample : I1182-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TOD-BC-0.2

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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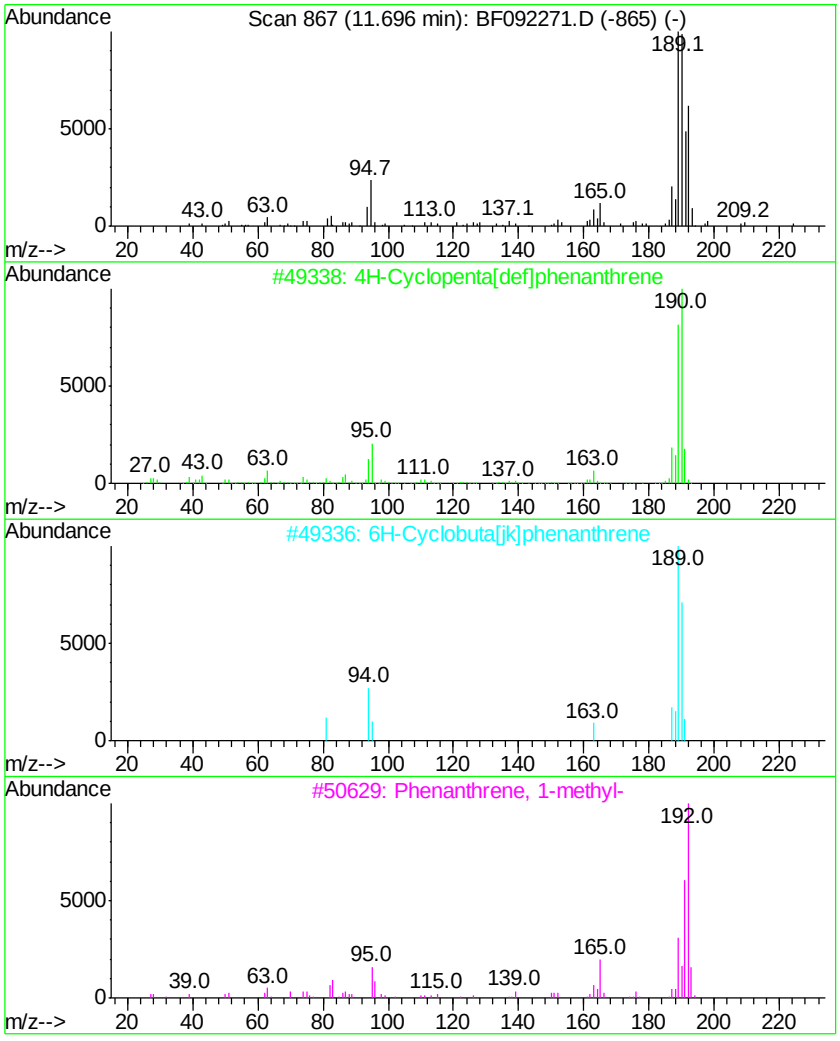
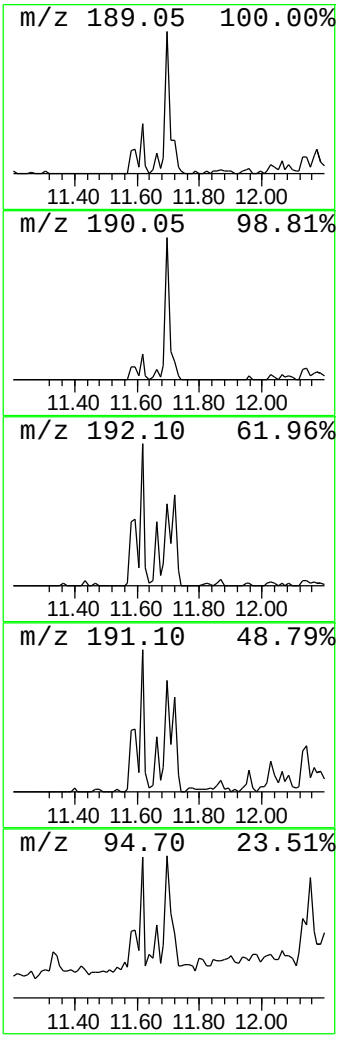
Peak Number 8 4H-Cyclopenta[def]phenanthrene Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.70	2.69 ng	172326	Phenanthrene-d10	11.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
2	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	40
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	38
4	3,5-Dichlorobenzoic acid	190	C7H4Cl2O2	000051-36-5	30
5	Naphtho[1,2-b]norbornadiene	192	C15H12	1000210-14-8	18





Data Path : Z:\HPCHEM1\BNA F\DATA\BF011317\
Data File : BF092271.D
Acq On : 14 Jan 2017 4:58
Operator : UM/SJ
Sample : I1182-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TOD-BC-0.2

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

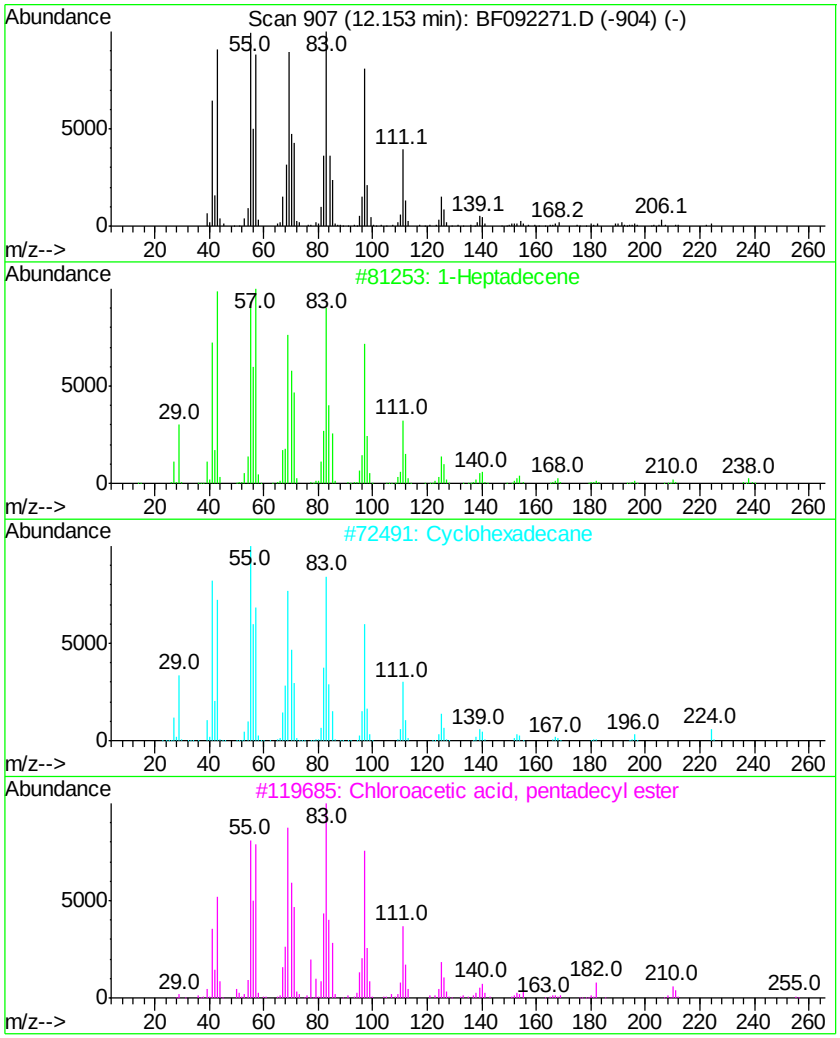
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

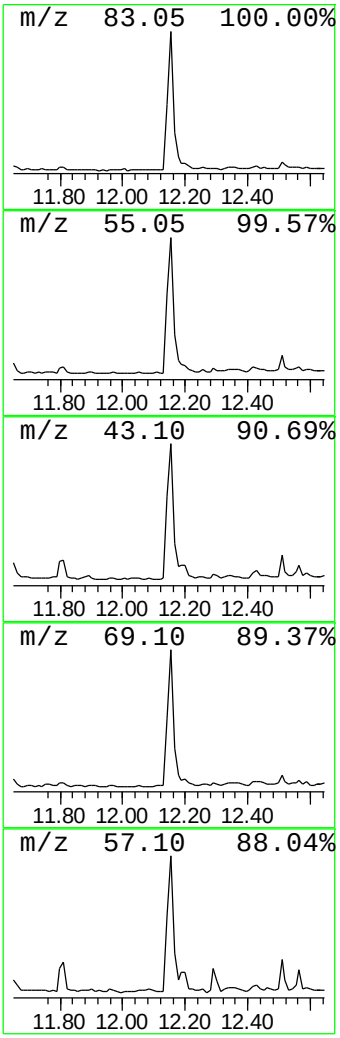
****

Peak Number 9 1-Heptadecene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.15	13.30 ng	853122	Phenanthrene-d10	11.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heptadecene	238	C17H34	006765-39-5	94
2	Cyclohexadecane	224	C16H32	000295-65-8	94
3	Chloroacetic acid, pentadecyl ester	304	C17H33ClO2	070301-47-2	94
4	1-Octadecanol	270	C18H38O	000112-92-5	94
5	1-Nonadecene	266	C19H38	018435-45-5	91





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011317\
Data File : BF092271.D
Acq On : 14 Jan 2017 4:58
Operator : UM/SJ
Sample : I1182-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TOD-BC-0.2

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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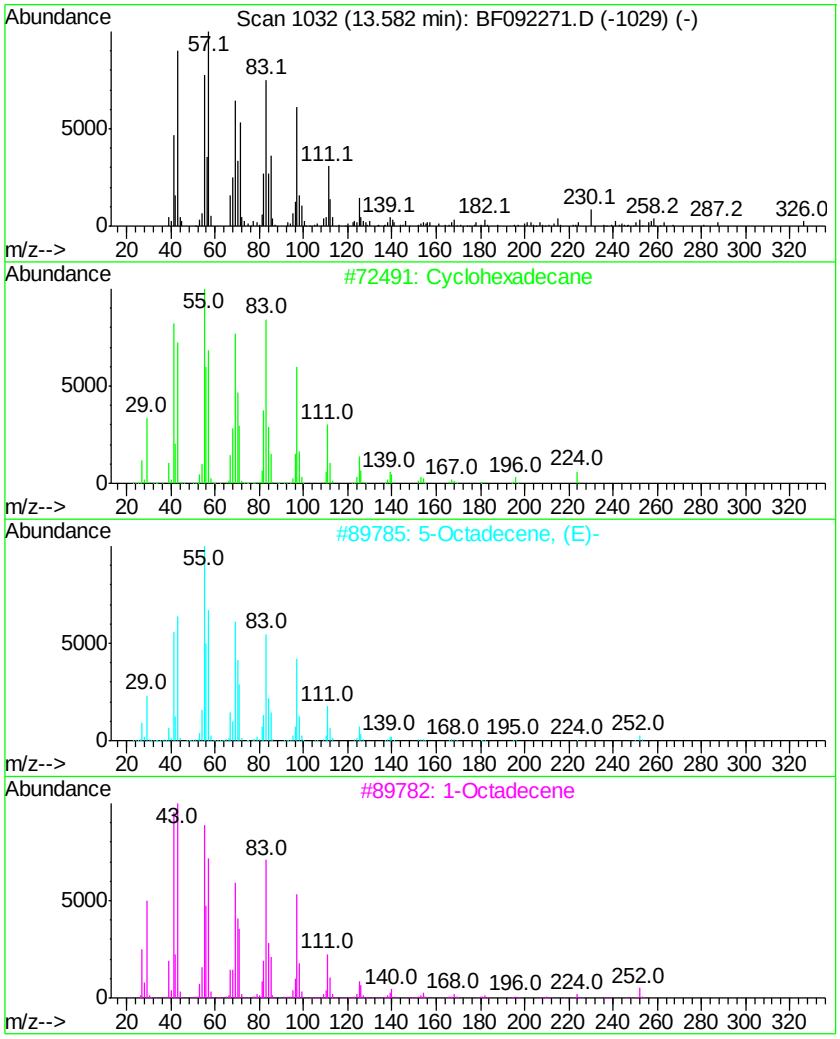
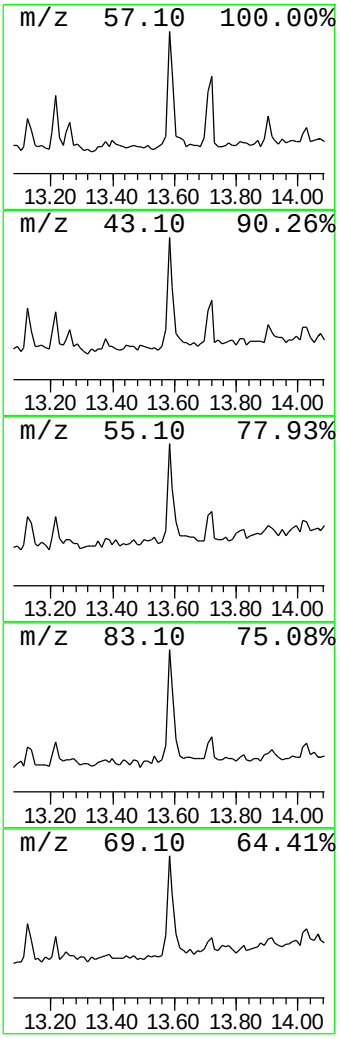
Peak Number 10 Cyclohexadecane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.58	2.79 ng	237423	Chrysene-d12	13.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexadecane	224	C16H32	000295-65-8	98
2	5-Octadecene, (E)-	252	C18H36	007206-21-5	95
3	1-Octadecene	252	C18H36	000112-88-9	95
4	Heptafluorobutanoic acid, heptad...	452	C21H35F7O2	1000282-97-3	94
5	1-Tricosene	322	C23H46	018835-32-0	91





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Data File : BF092271.D
Acq On : 14 Jan 2017 4:58
Operator : UM/SJ
Sample : I1182-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TOD-BC-0.2

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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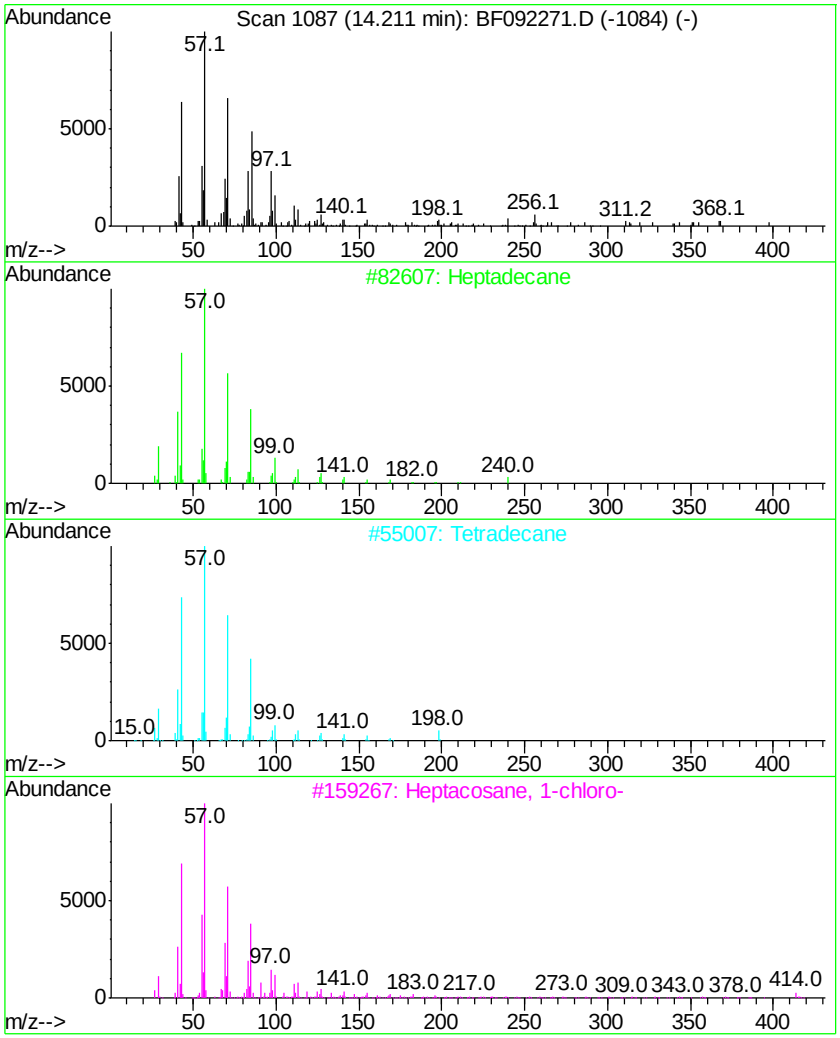
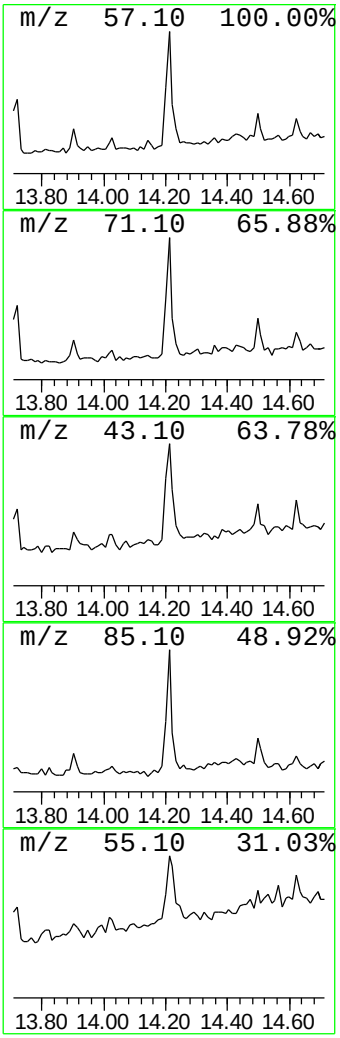
Peak Number 11 Heptadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.21	3.16 ng	269008	Chrysene-d12	13.72

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	96
2	Tetradecane	198	C14H30	000629-59-4	91
3	Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	80
4	Nonadecane, 9-methyl-	282	C20H42	013287-24-6	76
5	Tritetracontane	605	C43H88	007098-21-7	74





Library Search Compound Report

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Operator : UM/SJ
Sample : I1182-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TOD-BC-0.2

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
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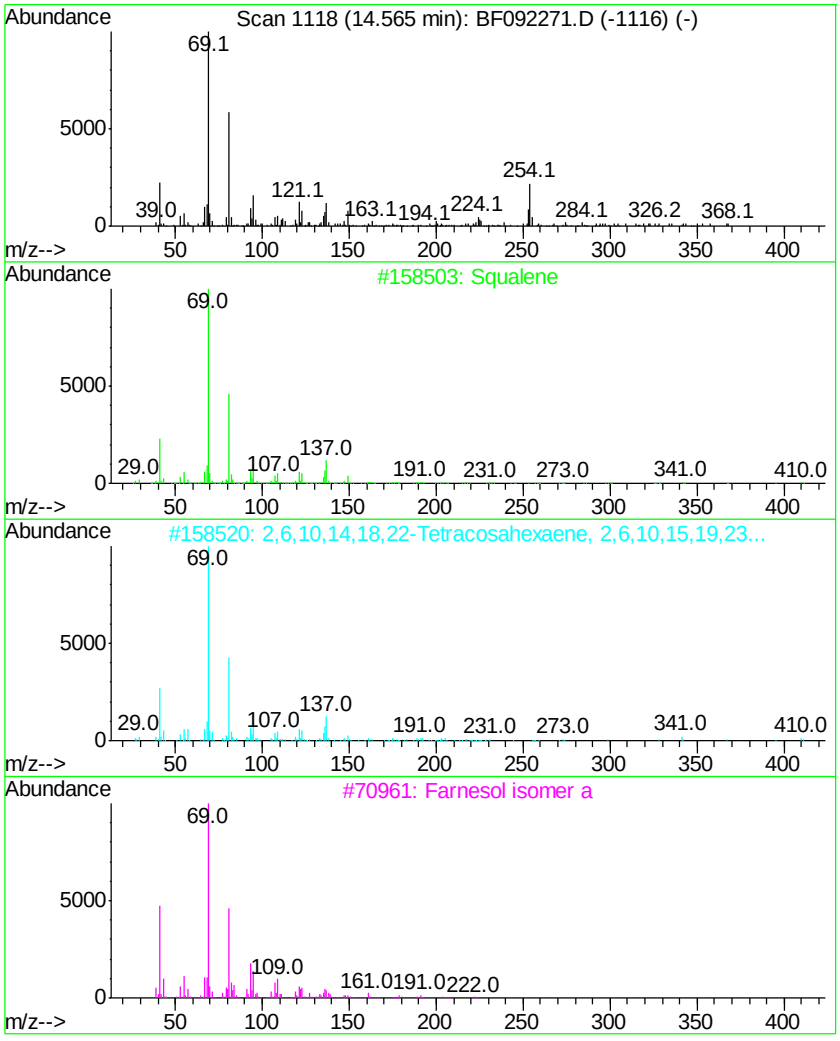
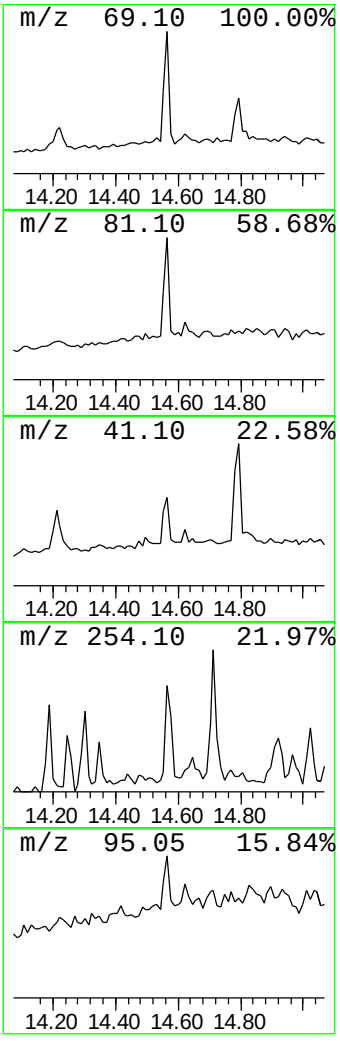
Peak Number 12 Squalene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.57	3.64 ng	151447	Perylene-d12	15.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Squalene	410	C30H50	007683-64-9	68
2	2,6,10,14,18,22-Tetracosahexaene...	410	C30H50	000111-02-4	64
3	Farnesol isomer a	222	C15H26O	1000108-92-4	53
4	7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	53
5	4,8,12-Tetradecatrienenitrile, 5...	245	C17H27N	005981-31-7	50





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Data File : BF092271.D
Acq On : 14 Jan 2017 4:58
Operator : UM/SJ
Sample : I1182-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TOD-BC-0.2

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
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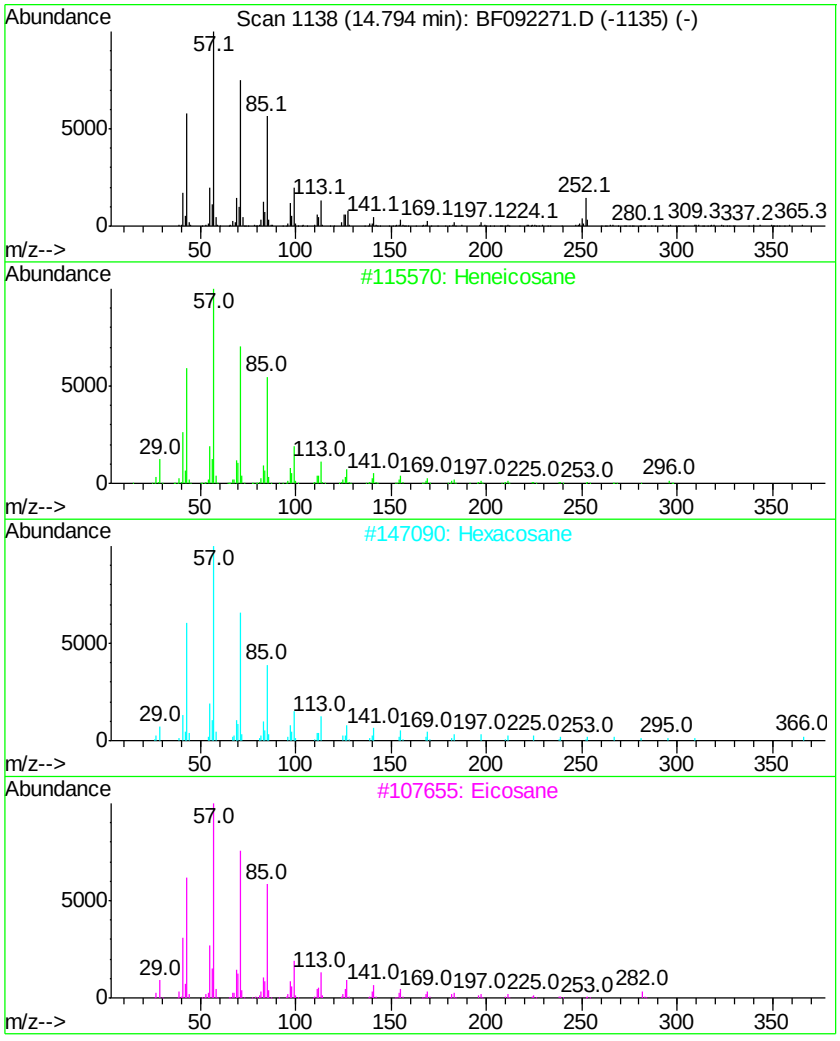
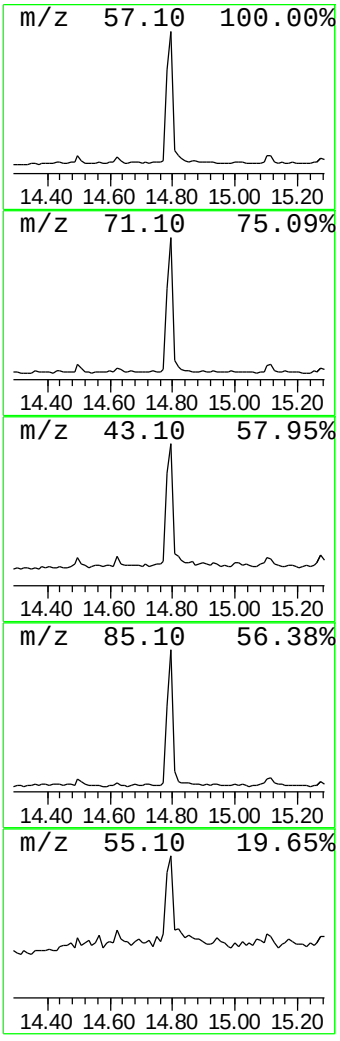
Peak Number 13 Heneicosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.79	11.86 ng	493763	Perylene-d12	15.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	90
2	Hexacosane	366	C26H54	000630-01-3	87
3	Eicosane	282	C20H42	000112-95-8	87
4	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	86
5	Pentacosane	352	C25H52	000629-99-2	86





Library Search Compound Report

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Sample : I1182-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TOD-BC-0.2

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
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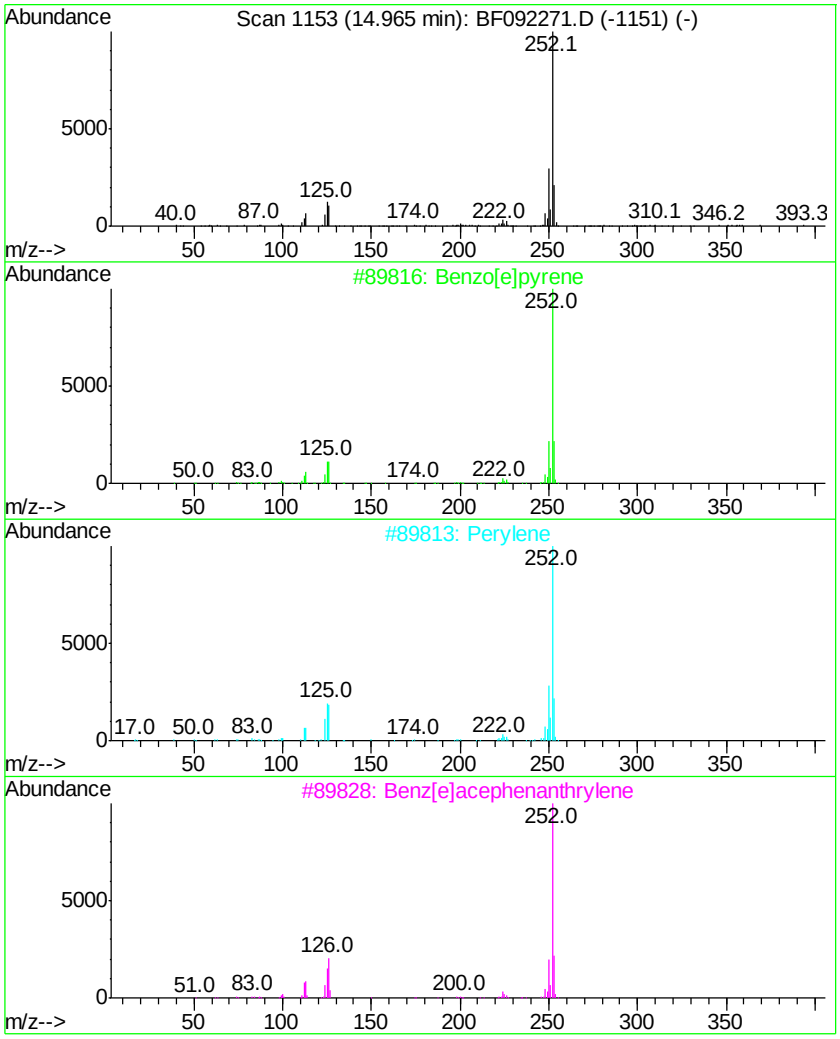
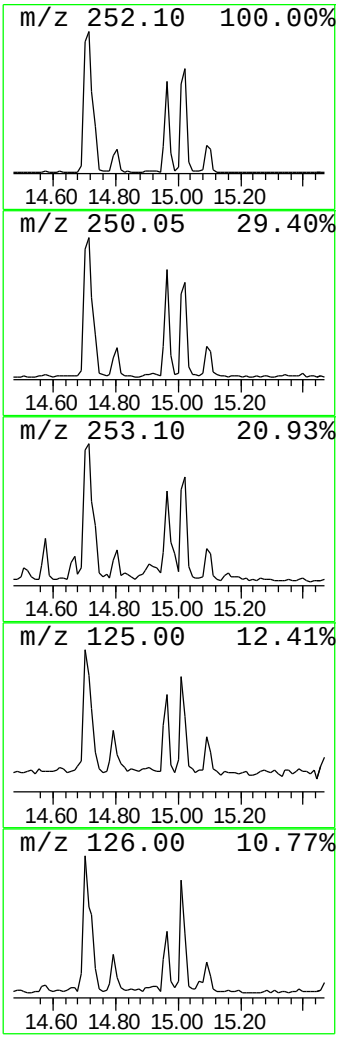
Peak Number 14 Benzo[e]pyrene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.97	3.70 ng	154229	Perylene-d12	15.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	99
2	Perylene	252	C20H12	000198-55-0	97
3	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	95
4	Benzo[a]pyrene	252	C20H12	000050-32-8	94
5	Benzo[k]fluoranthene	252	C20H12	000207-08-9	91





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ALS Vial : 32 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TOD-BC-0.2

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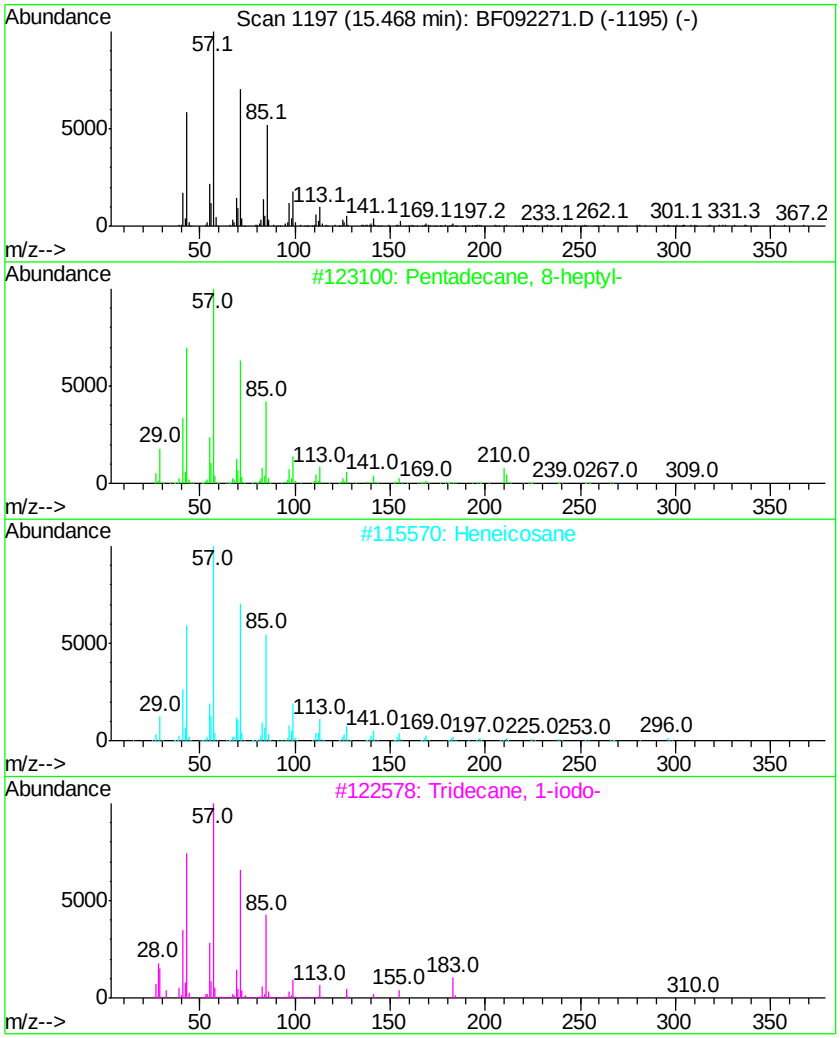
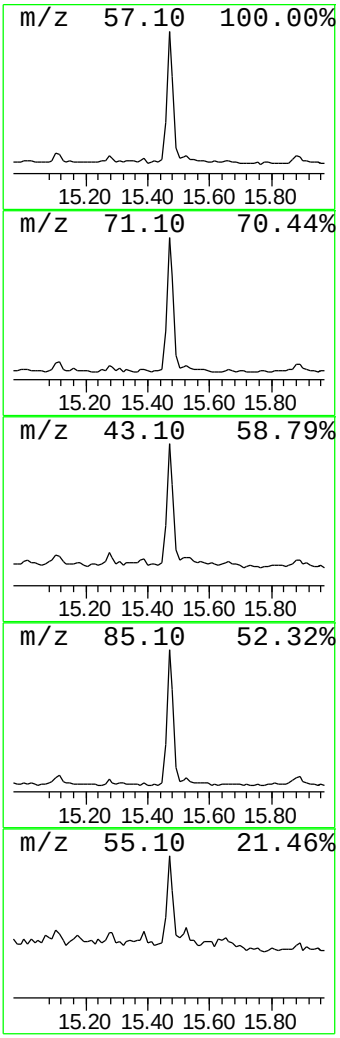
Peak Number 15 Pentadecane, 8-heptyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.47	11.30 ng	470671	Perylene-d12	15.07

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 8-heptyl-	310	C22H46	071005-15-7	91
2	Heneicosane	296	C21H44	000629-94-7	91
3	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	90
4	Hexadecane, 3-methyl-	240	C17H36	006418-43-5	87
5	Nonadecane	268	C19H40	000629-92-5	86





Tentatively Identified Compound (LSC) summary

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Acq On : 14 Jan 2017 4:58
Operator : UM/SJ
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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...	4.70	28.6	ng	1594540	1	6.58	1114330	20.0
unknown6.35	6.35	58.5	ng	3260910	1	6.58	1114330	20.0
Cyclotetradecane	11.33	5.6	ng	360698	4	11.09	1283200	20.0
Anthracene, 1-met...	11.58	2.0	ng	130973	4	11.09	1283200	20.0
n-Hexadecanoic acid	11.64	2.8	ng	180456	4	11.09	1283200	20.0
4H-Cyclopenta[def...	11.70	2.7	ng	172326	4	11.09	1283200	20.0
1-Heptadecene	12.15	13.3	ng	853122	4	11.09	1283200	20.0
Cyclohexadecane	13.58	2.8	ng	237423	5	13.72	1700490	20.0
Heptadecane	14.21	3.2	ng	269008	5	13.72	1700490	20.0
Squalene	14.57	3.6	ng	151447	6	15.07	832784	20.0
Heneicosane	14.79	11.9	ng	493763	6	15.07	832784	20.0
Benzo[e]pyrene	14.97	3.7	ng	154229	6	15.07	832784	20.0
Pentadecane, 8-he...	15.47	11.3	ng	470671	6	15.07	832784	20.0