LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032816\
Data File : BF085812.D
Acq On : 28 Mar 2016 20:16
Operator : UM/SJ
Sample : H1905-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TRACK-D-GRID-10C

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-BF032416.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	4.996	234	237	247	rBV	601297	919661	31.03%	3.241%
2	5.419	271	274	276	rBV	2354547	2563364	86.48%	9.034%
3	5.819	307	309	312	rBV	85289	77862	2.63%	0.274%
4	6.447	362	364	367	rBV	1730808	2593137	87.48%	9.139%
5	6.585	373	376	378	rBV	2759560	2677615	90.33%	9.436%
6	6.813	394	396	399	rBV	814085	689379	23.26%	2.429%
7	6.973	407	410	412	rBV	2249938	2211623	74.61%	7.794%
8	7.385	443	446	448	rBV	1674585	1850702	62.44%	6.522%
9	8.105	506	509	511	rBV	720598	865353	29.19%	3.050%
10	9.179	600	603	605	rBV	3170290	2964126	100.00%	10.446%
11	9.579	635	638	639	rBV	222280	292887	9.88%	1.032%
12	9.785	654	656	660	rBV	56512	66858	2.26%	0.236%
13	9.853	660	662	665	rVB	1093898	954502	32.20%	3.364%
14	10.288	696	700	702	rBV2	128811	122907	4.15%	0.433%
15	10.505	716	719	722	rBV	36505	60698	2.05%	0.214%
16	10.642	728	731	734	rBV	1609483	1708264	57.63%	6.020%
17	10.756	740	741	746	rVB	105005	138608	4.68%	0.488%
18	11.202	779	780	782	rBV	42804	57337	1.93%	0.202%
19	11.339	790	792	794	rBV	1102062	957954	32.32%	3.376%
20	11.568	810	812	816	rBV	29192	43905	1.48%	0.155%
21	11.865	837	838	845	rBV2	58405	83541	2.82%	0.294%
22	12.654	904	907	911	rBV2	23772	37302	1.26%	0.131%
23	12.745	913	915	917	rBV	35388	32757	1.11%	0.115%
24	12.928	928	931	933	rBV	2261189	2888549	97.45%	10.180%
25	13.157	948	951	953	rBV	25761	32312	1.09%	0.114%
26	13.397	970	972	978	rBV	122076	348879	11.77%	1.230%
27	13.488	978	980	984	rVB2	36293	68793	2.32%	0.242%
28	13.820	1006	1009	1018	rBV2	121475	183304	6.18%	0.646%
29	13.968	1020	1022	1025	rVB	592209	633900	21.39%	2.234%
30	14.140	1035	1037	1042	rVB	54400	61582	2.08%	0.217%
31	14.437	1061	1063	1066	rBV	40542	58571	1.98%	0.206%
32	14.734	1087	1089	1091	rBV	41094	46259	1.56%	0.163%
33	14.837	1095	1098	1100	rVB3	21925	50649	1.71%	0.178%
34	15.008	1108	1113	1115	rBV4	19921	63078	2.13%	0.222%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032816\
Data File : BF085812.D
Acq On : 28 Mar 2016 20:16
Operator : UM/SJ
Sample : H1905-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TRACK-D-GRID-10C

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-BF032416.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	15.054	1115	1117	1119	rVV	45230	65495	2.21%	0.231%
36	15.145	1122	1125	1127	rVB2	36075	47628	1.61%	0.168%
37	15.283	1133	1137	1140	rBV5	51639	108083	3.65%	0.381%
38	15.385	1144	1146	1150	rVV	381863	612918	20.68%	2.160%
39	15.454	1150	1152	1156	rVV5	57612	106790	3.60%	0.376%
40	15.523	1156	1158	1160	rVB3	52652	66081	2.23%	0.233%
41	15.923	1190	1193	1200	rVV7	19380	58281	1.97%	0.205%
42	16.025	1200	1202	1206	rVB5	20535	40051	1.35%	0.141%
43	16.288	1221	1225	1228	rBV3	18193	48139	1.62%	0.170%
44	16.506	1241	1244	1247	rBV4	17298	37014	1.25%	0.130%
45	16.894	1275	1278	1281	rVB5	32144	40242	1.36%	0.142%
46	17.237	1307	1308	1311	rBV3	42447	41559	1.40%	0.146%
47	18.060	1377	1380	1382	rBV4	17544	34514	1.16%	0.122%
48	18.231	1393	1395	1397	rBV2	25669	32464	1.10%	0.114%
49	18.300	1399	1401	1403	rVV2	24097	32987	1.11%	0.116%
50	18.346	1403	1405	1407	rVB3	34245	42468	1.43%	0.150%
51	18.883	1450	1452	1453	rBV2	39004	50709	1.71%	0.179%
52	18.906	1453	1454	1456	rVB2	39245	47680	1.61%	0.168%
53	19.157	1473	1476	1483	rVB4	61205	165327	5.58%	0.583%
54	19.272	1483	1486	1488	rBV3	51601	69418	2.34%	0.245%
55	19.443	1498	1501	1505	rBV4	53577	174584	5.89%	0.615%
56	19.523	1505	1508	1509	rBV2	21364	46855	1.58%	0.165%

Sum of corrected areas: 28375505

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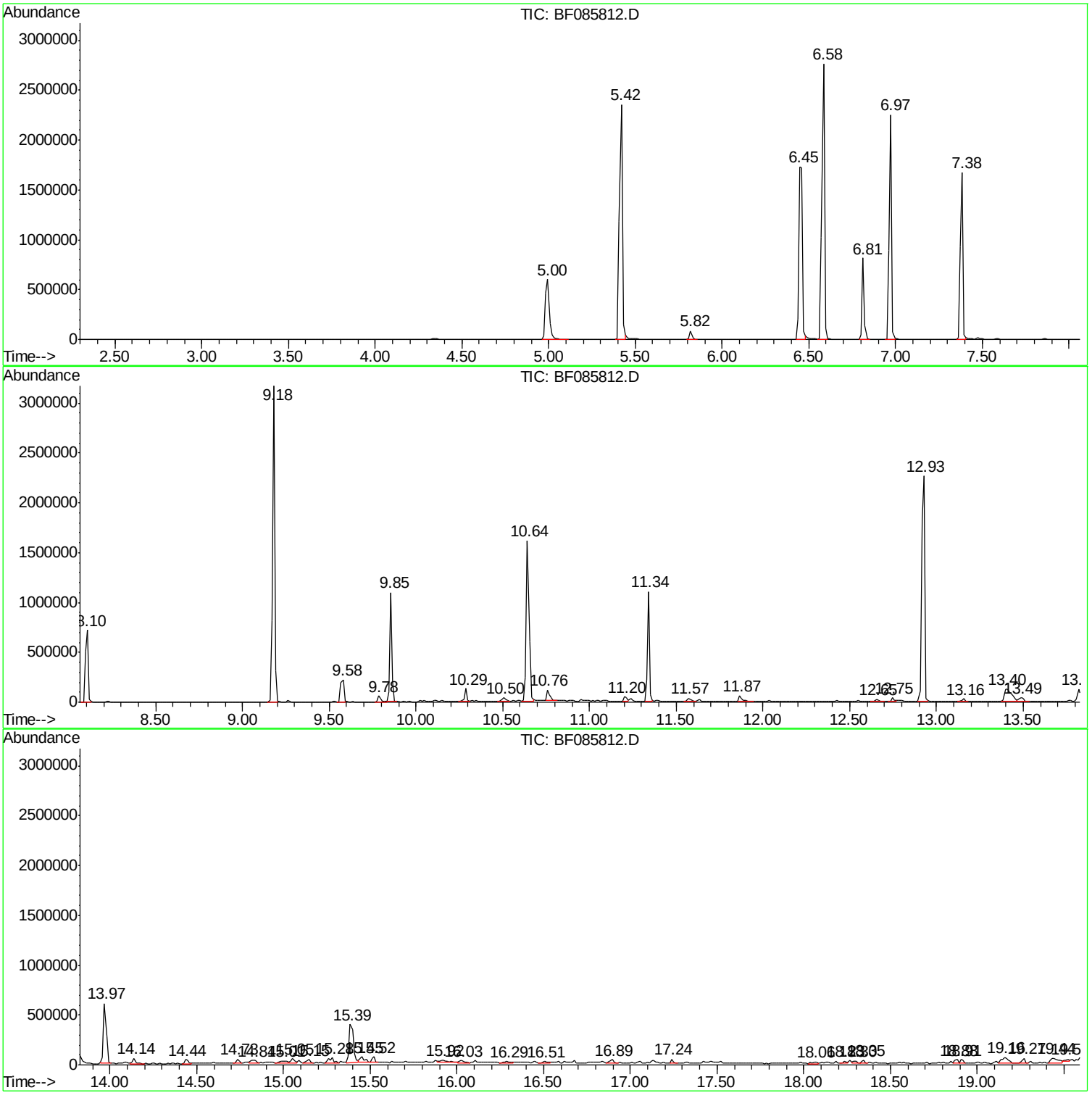
Data File : BF085812.D
Acq On : 28 Mar 2016 20:16
Operator : UM/SJ
Sample : H1905-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TRACK-D-GRID-10C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.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\BF032816\
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Acq On : 28 Mar 2016 20:16
Operator : UM/SJ
Sample : H1905-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TRACK-D-GRID-10C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

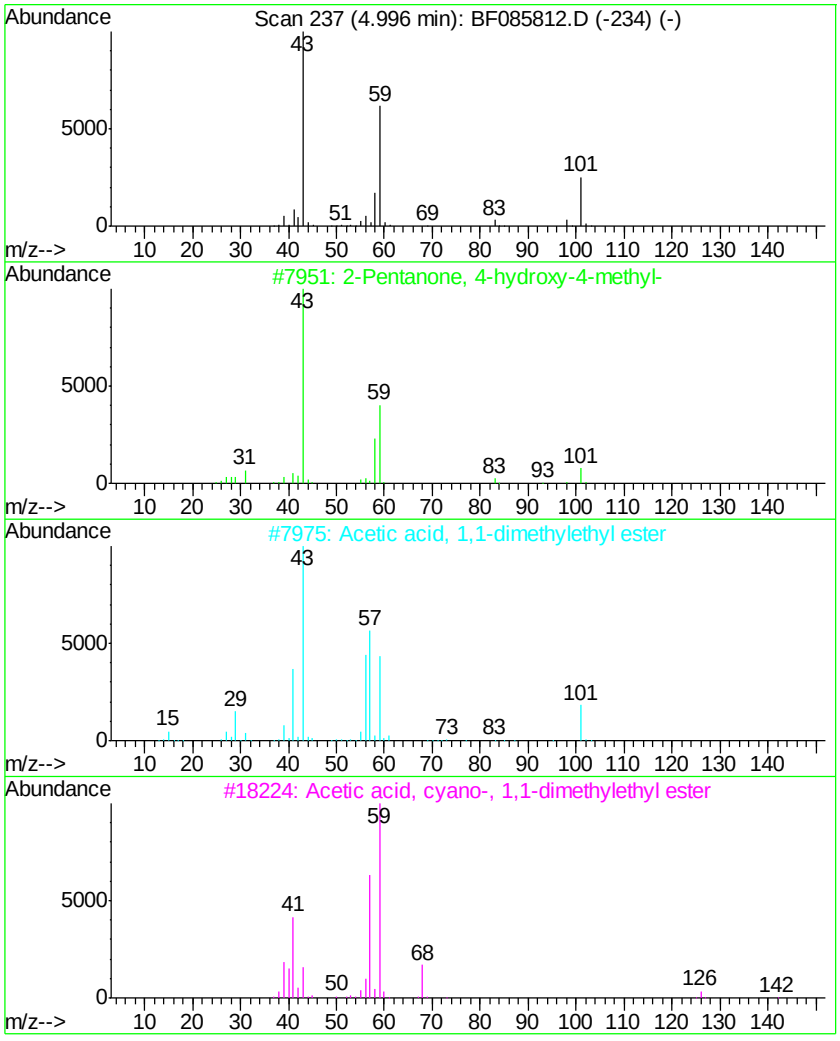
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

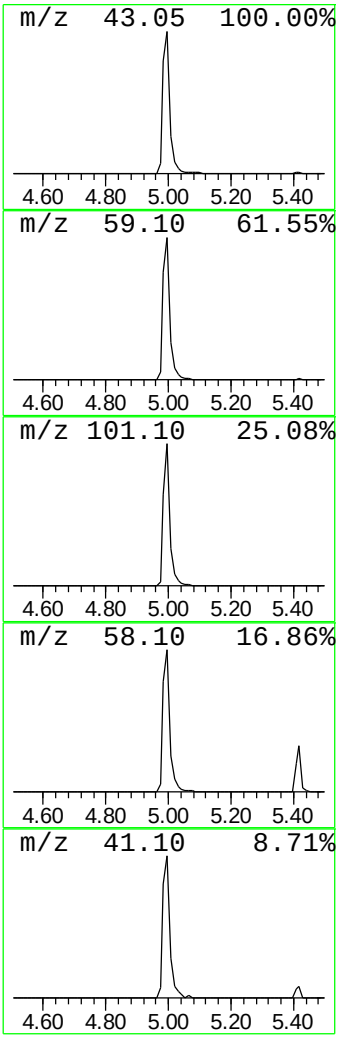
****

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.00	26.68 ng	919661	1,4-Dichlorobenzene-d4	6.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
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	5-Hexen-2-one	98	C6H10O	000109-49-9	9
5	(+-)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032816\
Data File : BF085812.D
Acq On : 28 Mar 2016 20:16
Operator : UM/SJ
Sample : H1905-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TRACK-D-GRID-10C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.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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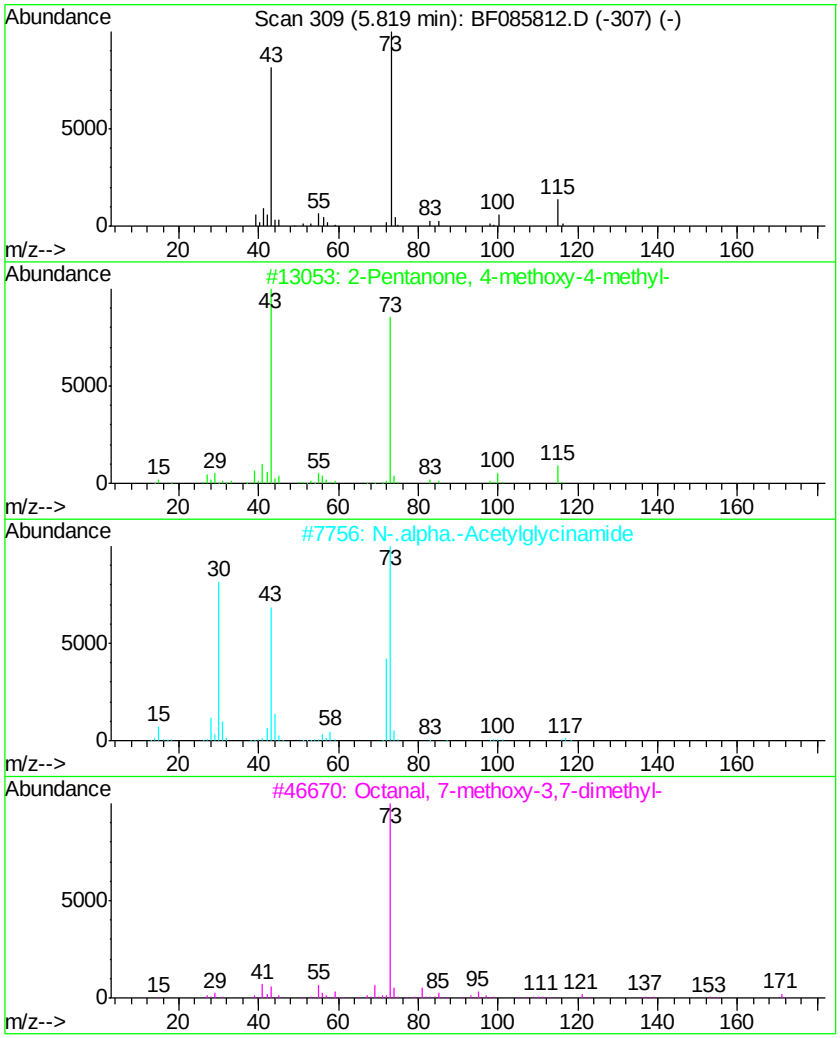
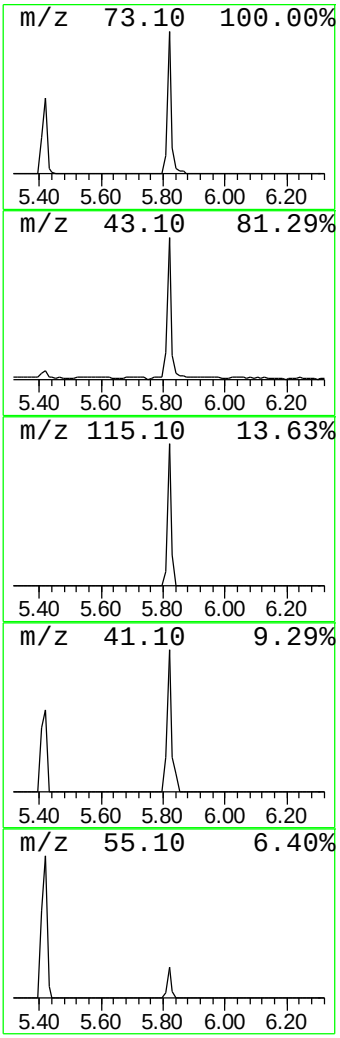
Peak Number 2 2-Pentanone, 4-methoxy-4-me... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.82	2.26 ng	77862	1,4-Dichlorobenzene-d4	6.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-methoxy-4-methyl-	130	C7H14O2	000107-70-0	83
2	N-.alpha.-Acetylglycinamide	116	C4H8N2O2	002620-63-5	33
3	Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	17
4	Pentanoic acid, 4-oxo-	116	C5H8O3	000123-76-2	10
5	2-Pentanone, 3-methyl-	100	C6H12O	000565-61-7	10





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032816\
Data File : BF085812.D
Acq On : 28 Mar 2016 20:16
Operator : UM/SJ
Sample : H1905-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TRACK-D-GRID-10C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.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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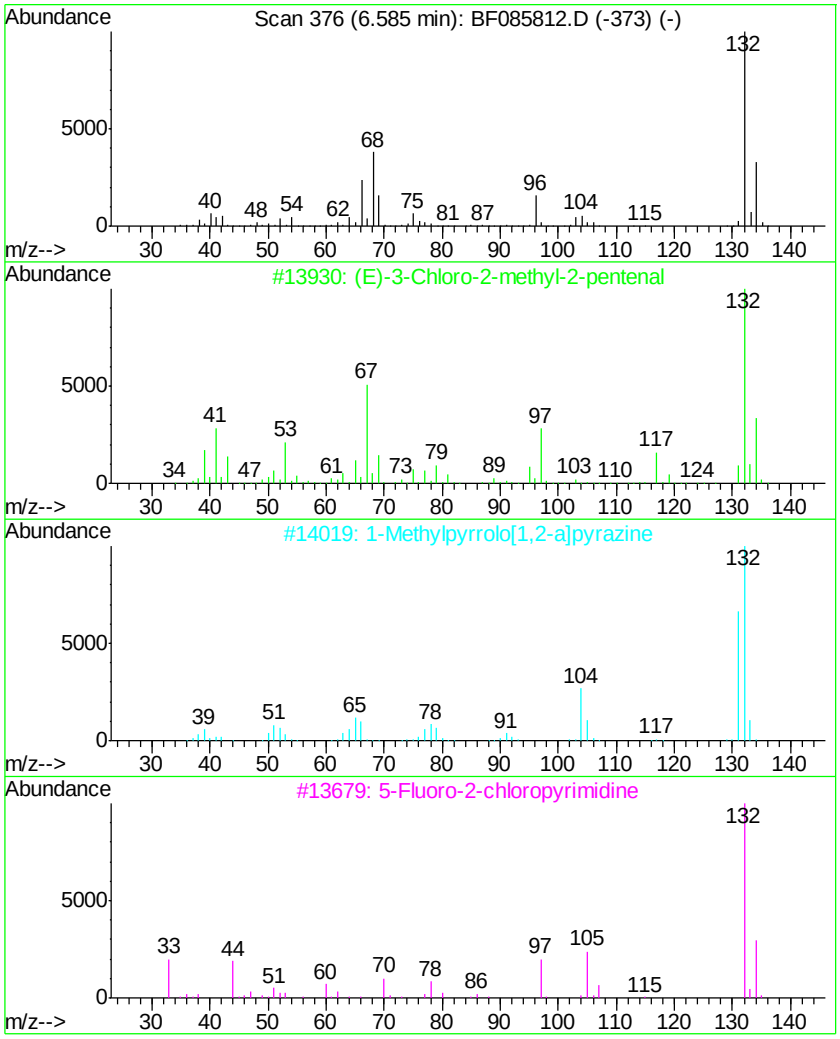
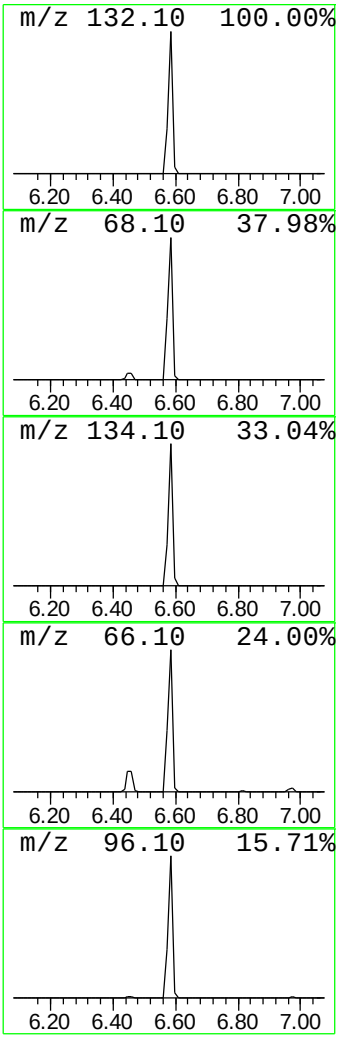
Peak Number 3 unknown6.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	77.68 ng	2677620	1,4-Dichlorobenzene-d4	6.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
2	1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
3	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
4	2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	9





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032816\
Data File : BF085812.D
Acq On : 28 Mar 2016 20:16
Operator : UM/SJ
Sample : H1905-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TRACK-D-GRID-10C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.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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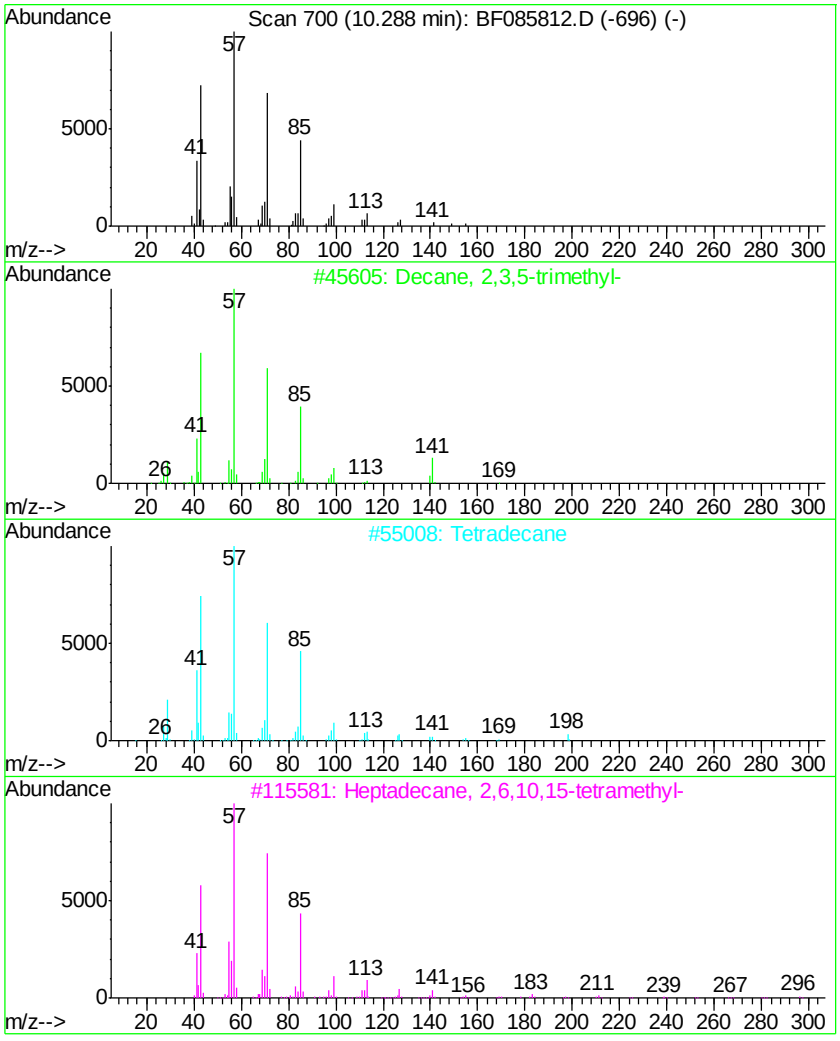
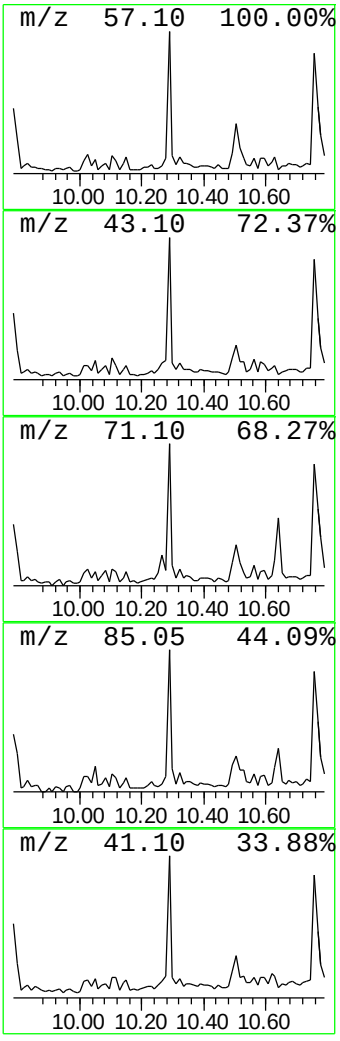
Peak Number 5 Decane, 2,3,5-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.29	2.58 ng	122907	Acenaphthene-d10	9.85

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	90
2	Tetradecane	198	C14H30	000629-59-4	87
3	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	87
4	Heneicosane	296	C21H44	000629-94-7	87
5	Nonadecane	268	C19H40	000629-92-5	86





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Operator : UM/SJ
Sample : H1905-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TRACK-D-GRID-10C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.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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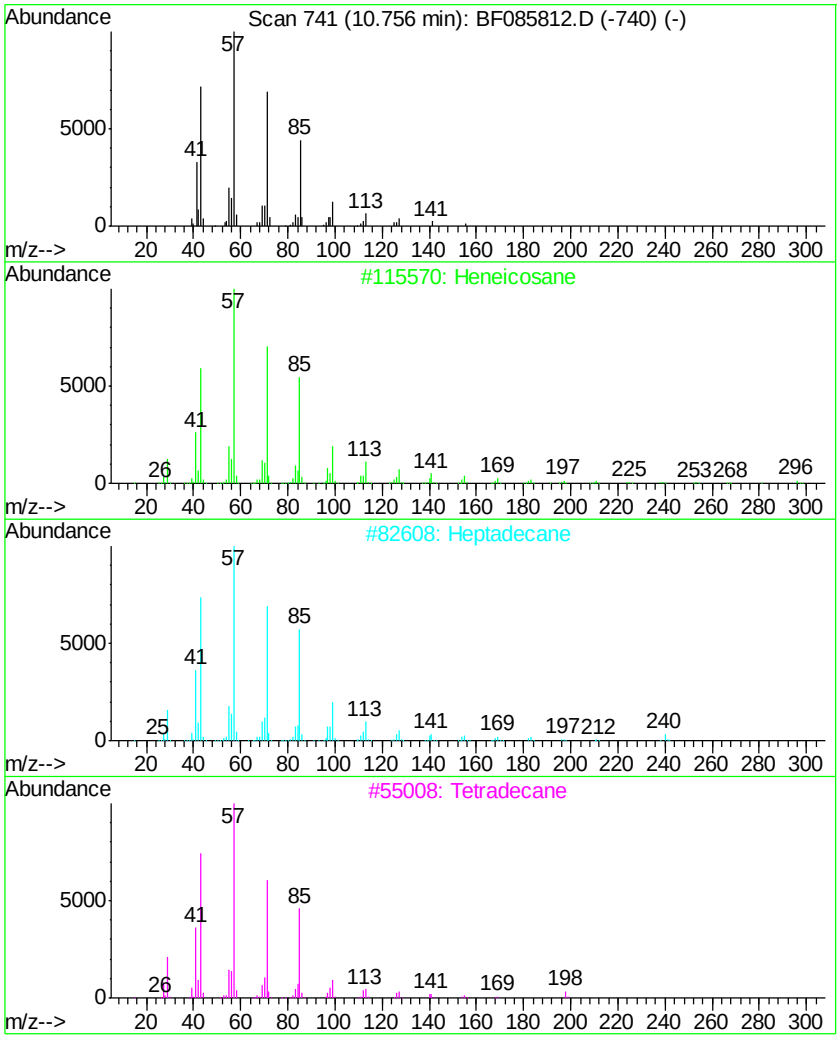
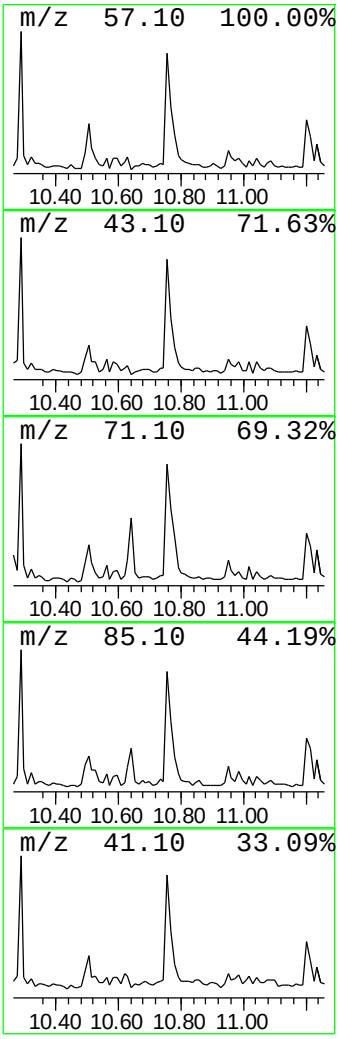
Peak Number 6 Heneicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.76	2.89 ng	138608	Phenanthrene-d10	11.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	90
2	Heptadecane	240	C17H36	000629-78-7	87
3	Tetradecane	198	C14H30	000629-59-4	86
4	Eicosane	282	C20H42	000112-95-8	86
5	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86





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Operator : UM/SJ
Sample : H1905-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TRACK-D-GRID-10C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.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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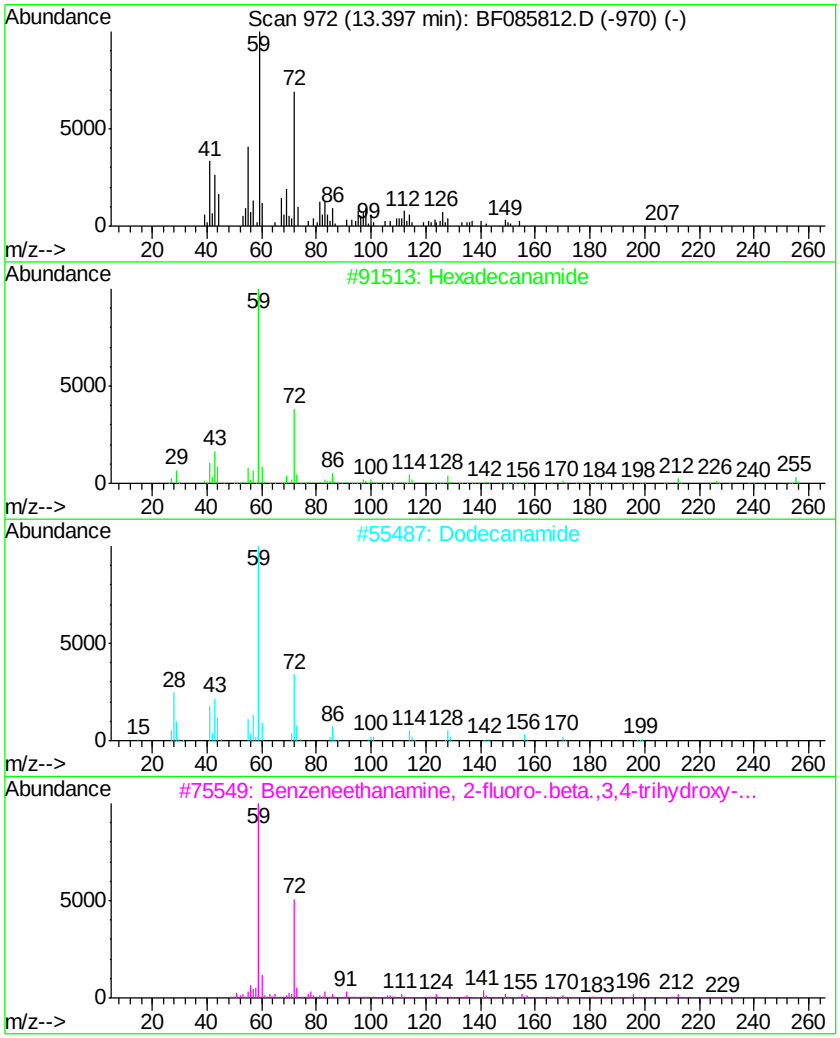
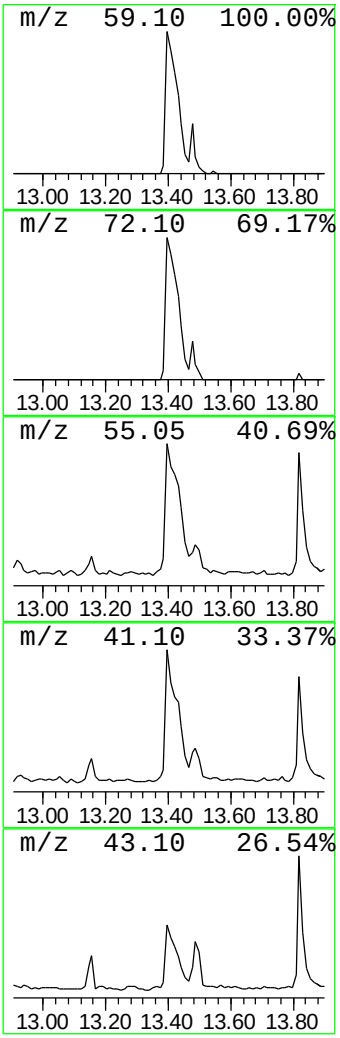
Peak Number 7 Hexadecanamide Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.40	11.01 ng	348879	Chrysene-d12	13.97

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecanamide	255	C16H33NO	000629-54-9	72
2	Dodecanamide	199	C12H25NO	001120-16-7	59
3	Benzeneethanamine, 2-fluoro-.bet...	229	C11H16FNO3	061338-98-5	59
4	Nonanamide	157	C9H19NO	001120-07-6	59
5	Octanamide	143	C8H17NO	000629-01-6	59





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Acq On : 28 Mar 2016 20:16
Operator : UM/SJ
Sample : H1905-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TRACK-D-GRID-10C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.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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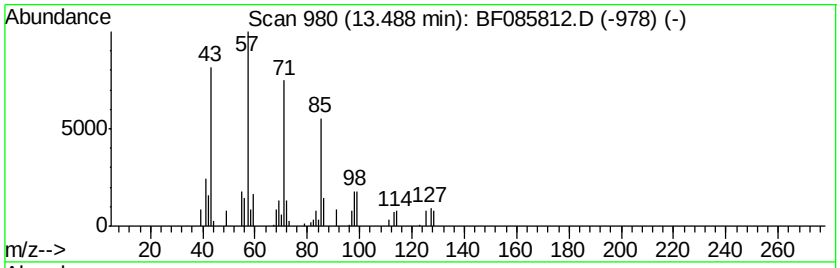
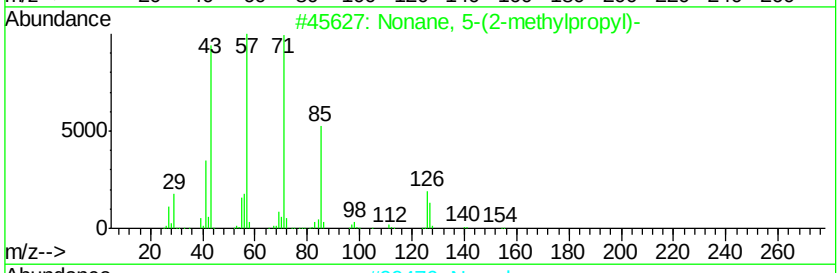
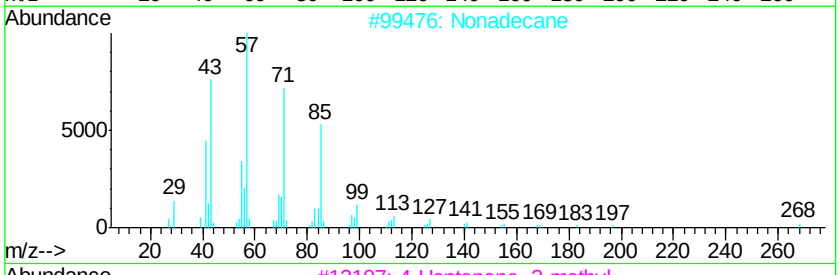
Peak Number 8 Nonane, 5-(2-methylpropyl)- Concentration Rank 14

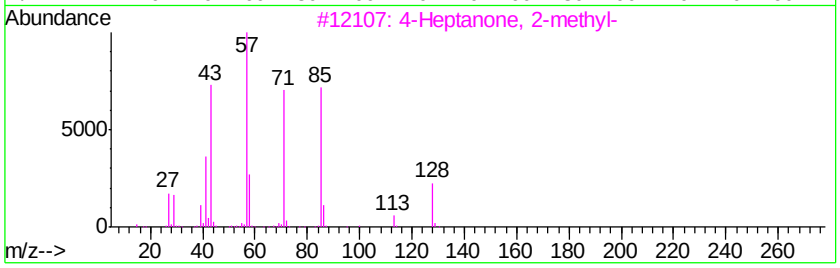
R.T.	EstConc	Area	Relative to ISTD	R.T.
13.49	2.17 ng	68793	Chrysene-d12	13.97

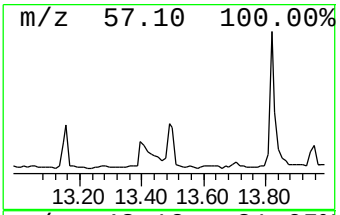
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane, 5-(2-methylpropyl)-	184	C13H28	062185-53-9	59
2	Nonadecane	268	C19H40	000629-92-5	59
3	4-Heptanone, 2-methyl-	128	C8H16O	000626-33-5	59
4	3-Octen-2-ol, 2-methyl-, (Z)-	142	C9H18O	018521-07-8	50
5	Eicosane	282	C20H42	000112-95-8	50

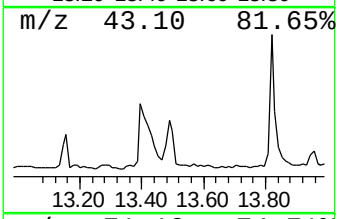


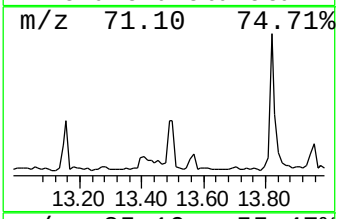


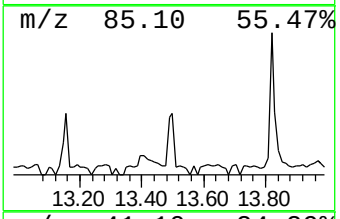


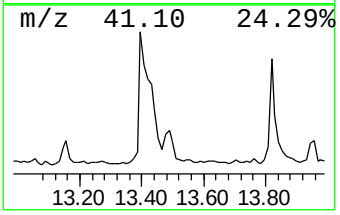












Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032816\
Data File : BF085812.D
Acq On : 28 Mar 2016 20:16
Operator : UM/SJ
Sample : H1905-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TRACK-D-GRID-10C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.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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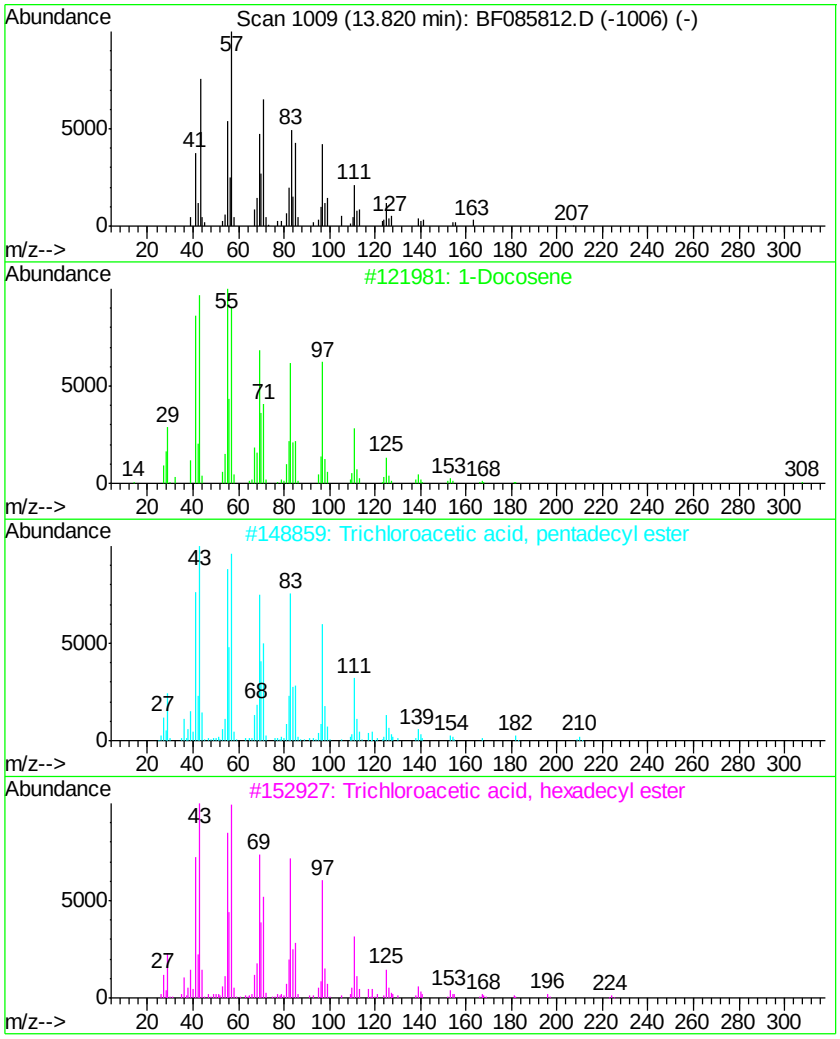
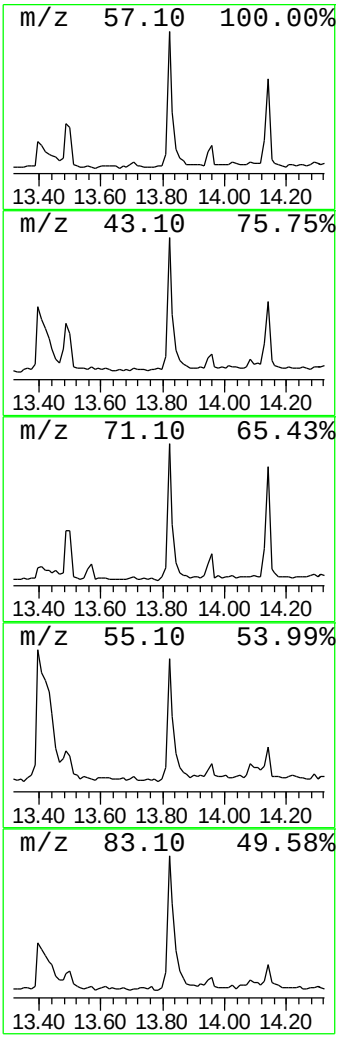
Peak Number 9 1-Docosene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.82	5.78 ng	183304	Chrysene-d12	13.97

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	76
2	Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	70
3	Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	70
4	3-Chloropropionic acid, heptadec...	346	C20H39ClO2	1000283-05-1	70
5	17-Pentatriacontene	491	C35H70	006971-40-0	68





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Acq On : 28 Mar 2016 20:16
Operator : UM/SJ
Sample : H1905-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TRACK-D-GRID-10C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.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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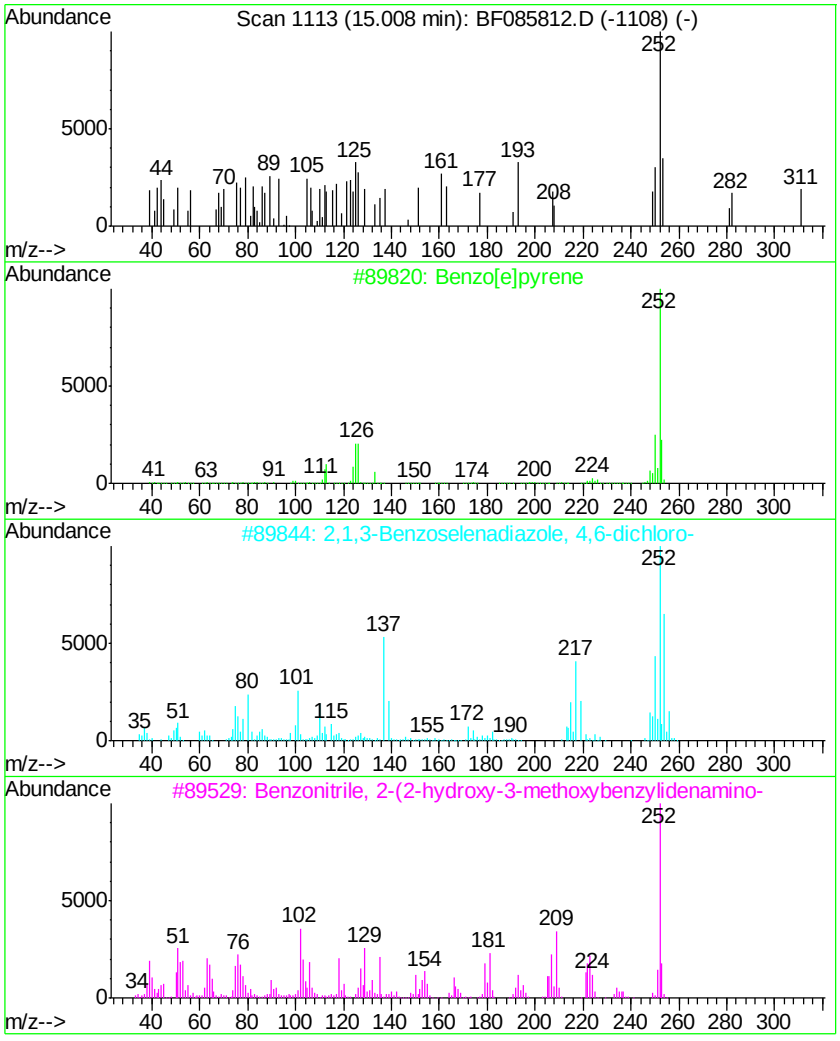
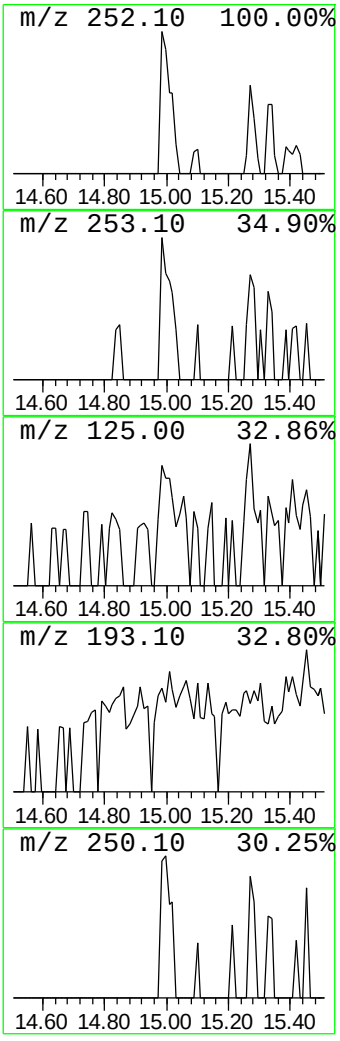
Peak Number 10 unknown15.01 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.01	2.06 ng	63078	Perylene-d12	15.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	38
2	2,1,3-Benzoselenadiazole, 4,6-di...	252	C6H2Cl2N2Se	090030-99-2	32
3	Benzonitrile, 2-(2-hydroxy-3-met...	252	C15H12N2O2	315670-28-1	27
4	Butanal, (2,4-dinitrophenyl)hydr...	252	C10H12N4O4	001527-98-6	27
5	1,2,4-Oxadiazole, 3,5-bis(4-meth...	282	C8H6N6O6	1000271-31-7	25





Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032816\
Data File : BF085812.D
Acq On : 28 Mar 2016 20:16
Operator : UM/SJ
Sample : H1905-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TRACK-D-GRID-10C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.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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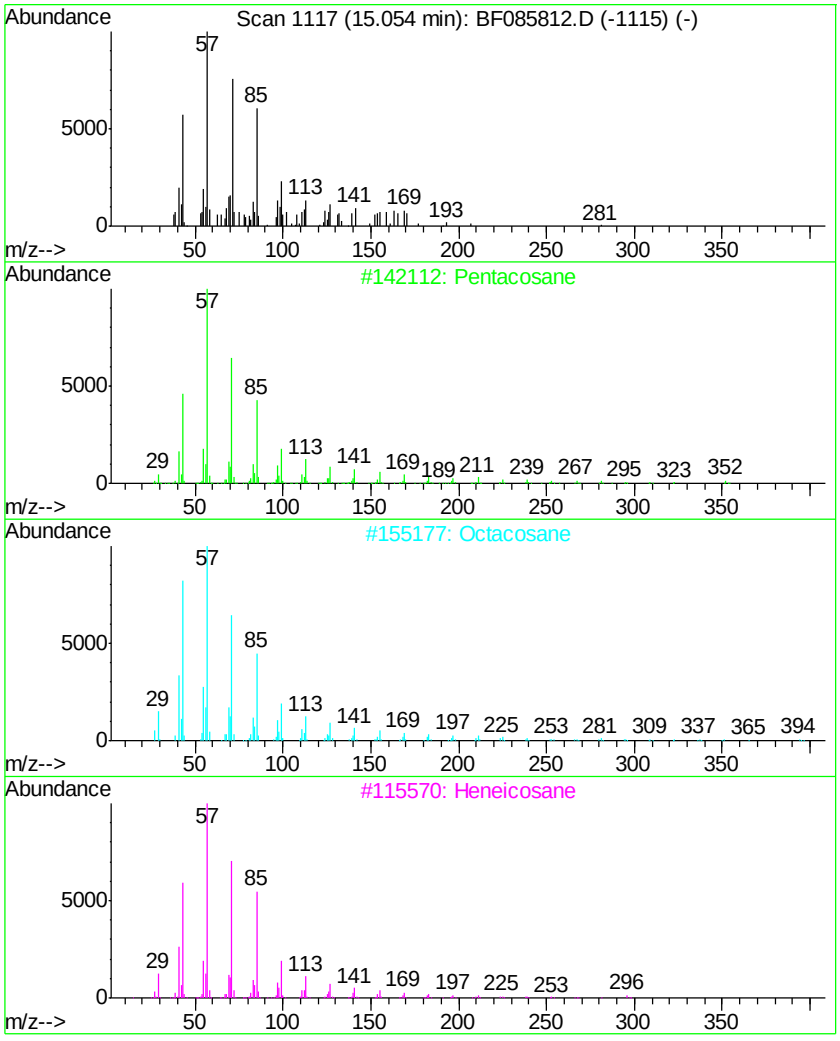
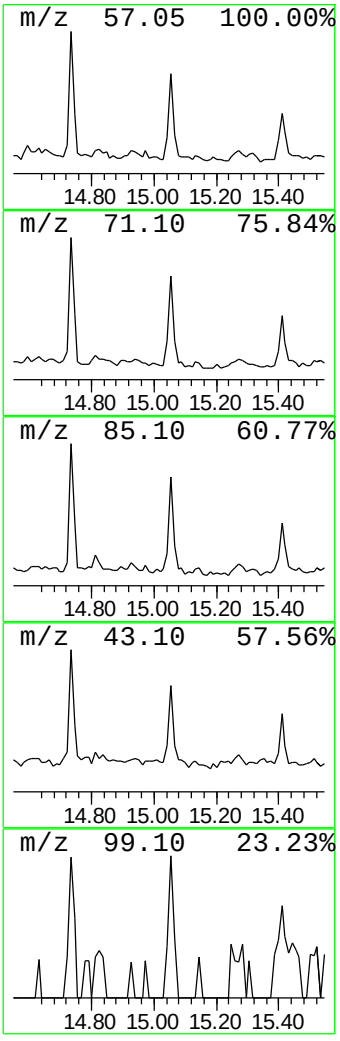
Peak Number 11 Pentacosane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.05	2.14 ng	65495	Perylene-d12	15.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentacosane	352	C25H52	000629-99-2	91
2	Octacosane	394	C28H58	000630-02-4	91
3	Heneicosane	296	C21H44	000629-94-7	91
4	Triacontane	422	C30H62	000638-68-6	90
5	Eicosane	282	C20H42	000112-95-8	90





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Data File : BF085812.D
Acq On : 28 Mar 2016 20:16
Operator : UM/SJ
Sample : H1905-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TRACK-D-GRID-10C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.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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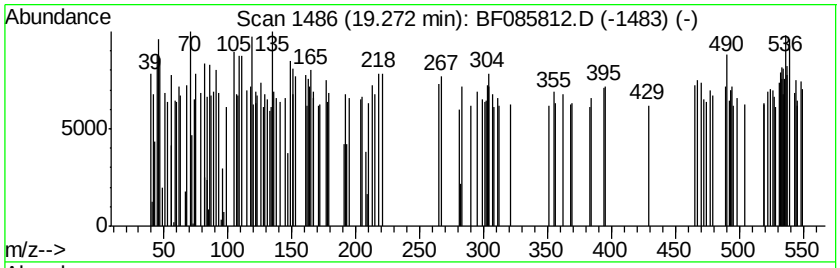
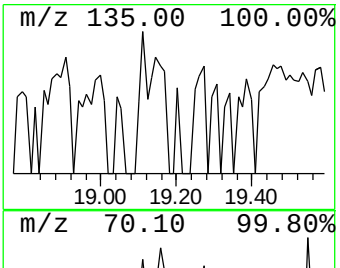
Peak Number 16 Iminobis[[bis(trimethylsily... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.27	2.27 ng	69418	Perylene-d12	15.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Iminobis[[bis(trimethylsilyl)ami...	533	C18H57B2N5Si6	113678-60-7	64
2	Osmium methyl, (.eta.-5-cyclopen...	548	C22H30OsP2	1000158-92-4	22
3	Phosphine, 4,6-dibenzofurandiylb...	536	C36H26OP2	133850-81-4	10
4	1,3,5,7,9,11,13-Heptaethyl-11-me...	562	C15H42O9Si7	073113-20-9	10
5	1,1',2,2',8,8',15,15'-Octamethyl...	536	C41H44	1000202-07-2	9





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032816\
Data File : BF085812.D
Acq On : 28 Mar 2016 20:16
Operator : UM/SJ
Sample : H1905-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TRACK-D-GRID-10C

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.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...	5.00	26.7	ng	919661	1	6.81	689379	20.0
2-Pentanone, 4-me...	5.82	2.3	ng	77862	1	6.81	689379	20.0
unknown6.58	6.58	77.7	ng	2677620	1	6.81	689379	20.0
Decane, 2,3,5-tri...	10.29	2.6	ng	122907	3	9.85	954502	20.0
Heneicosane	10.76	2.9	ng	138608	4	11.34	957954	20.0
Hexadecanamide	13.40	11.0	ng	348879	5	13.97	633900	20.0
Nonane, 5-(2-meth...	13.49	2.2	ng	68793	5	13.97	633900	20.0
1-Docosene	13.82	5.8	ng	183304	5	13.97	633900	20.0
unknown15.01	15.01	2.1	ng	63078	6	15.39	612918	20.0
Pentacosane	15.05	2.1	ng	65495	6	15.39	612918	20.0
Iminobis[[bis(tri...	19.27	2.3	ng	69418	6	15.39	612918	20.0