LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032816\
Data File : BF085805.D
Acq On : 28 Mar 2016 16:57
Operator : UM/SJ
Sample : H1941-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-8

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	4.333	178	179	183	rBV	24887	35579	1.16%	0.119%
2	4.996	234	237	245	rBV	964225	1415887	46.32%	4.721%
3	5.418	271	274	277	rBV	2562494	2634701	86.20%	8.785%
4	5.819	307	309	312	rBV	154862	134993	4.42%	0.450%
5	6.367	355	357	362	rBV	50481	65549	2.14%	0.219%
6	6.459	362	365	367	rBV	2594363	3056577	100.00%	10.191%
7	6.584	373	376	378	rBV	2947976	2813165	92.04%	9.380%
8	6.813	394	396	399	rBV	757507	650991	21.30%	2.171%
9	6.973	407	410	412	rBV	2584422	2386763	78.09%	7.958%
10	7.384	443	446	448	rBV	2015998	2043554	66.86%	6.814%
11	7.476	452	454	461	rVB	21125	35618	1.17%	0.119%
12	8.104	506	509	511	rBV	765370	838933	27.45%	2.797%
13	9.179	600	603	606	rBV	3111348	2871761	93.95%	9.575%
14	9.579	635	638	639	rBV	499420	513572	16.80%	1.712%
15	9.785	652	656	659	rBV2	83118	115064	3.76%	0.384%
16	9.853	660	662	665	rVB	1019166	914555	29.92%	3.049%
17	10.025	673	677	678	rBV	20695	34341	1.12%	0.115%
18	10.288	698	700	702	rVB	147470	116465	3.81%	0.388%
19	10.505	717	719	723	rBV	38998	64055	2.10%	0.214%
20	10.642	728	731	734	rBV2	1617864	1913998	62.62%	6.382%
21	10.756	740	741	746	rBV	106215	144808	4.74%	0.483%
22	11.202	779	780	782	rBV	37160	50956	1.67%	0.170%
23	11.339	790	792	794	rBV	1039312	892672	29.20%	2.976%
24	11.865	837	838	841	rBV	114007	178021	5.82%	0.594%
25	12.425	883	887	889	rBV	55866	63627	2.08%	0.212%
26	12.573	896	900	904	rBV3	16467	32804	1.07%	0.109%
27	12.653	904	907	913	rVV2	39785	67036	2.19%	0.224%
28	12.928	928	931	933	rBV	2440814	2789800	91.27%	9.302%
29	13.294	961	963	968	rVB	474627	449228	14.70%	1.498%
30	13.659	992	995	998	rBV4	13889	32892	1.08%	0.110%
31	13.716	998	1000	1002	rBV	96677	97114	3.18%	0.324%
32	13.819	1006	1009	1017	rVB	198090	251719	8.24%	0.839%
33	13.968	1020	1022	1025	rVB	526418	615132	20.12%	2.051%
34	14.082	1031	1032	1034	rBV	27846	32514	1.06%	0.108%

LSC Area Percent Report

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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	14.128	1034	1036	1041	rVB2	42222	75070	2.46%	0.250%
36	14.311	1049	1052	1060	rVB3	55223	92932	3.04%	0.310%
37	14.459	1060	1065	1070	rBV4	25057	56533	1.85%	0.188%
38	14.814	1094	1096	1098	rVB	40604	56685	1.85%	0.189%
39	15.122	1118	1123	1128	rVB6	32293	101079	3.31%	0.337%
40	15.385	1144	1146	1149	rBV	344728	506208	16.56%	1.688%
41	15.797	1180	1182	1186	rVB2	15783	30751	1.01%	0.103%
42	15.877	1186	1189	1192	rVV3	14631	42553	1.39%	0.142%
43	15.968	1194	1197	1200	rVV2	92476	157463	5.15%	0.525%
44	16.071	1204	1206	1216	rVB6	26935	72905	2.39%	0.243%
45	16.380	1230	1233	1236	rVB	30551	55550	1.82%	0.185%
46	17.305	1310	1314	1320	rBV6	46471	161815	5.29%	0.540%
47	17.968	1367	1372	1380	rVB	49361	127324	4.17%	0.425%
48	18.254	1393	1397	1403	rBV3	22573	55894	1.83%	0.186%
49	18.368	1403	1407	1415	rVB6	11926	44370	1.45%	0.148%

Sum of corrected areas: 29991576

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BNA_F
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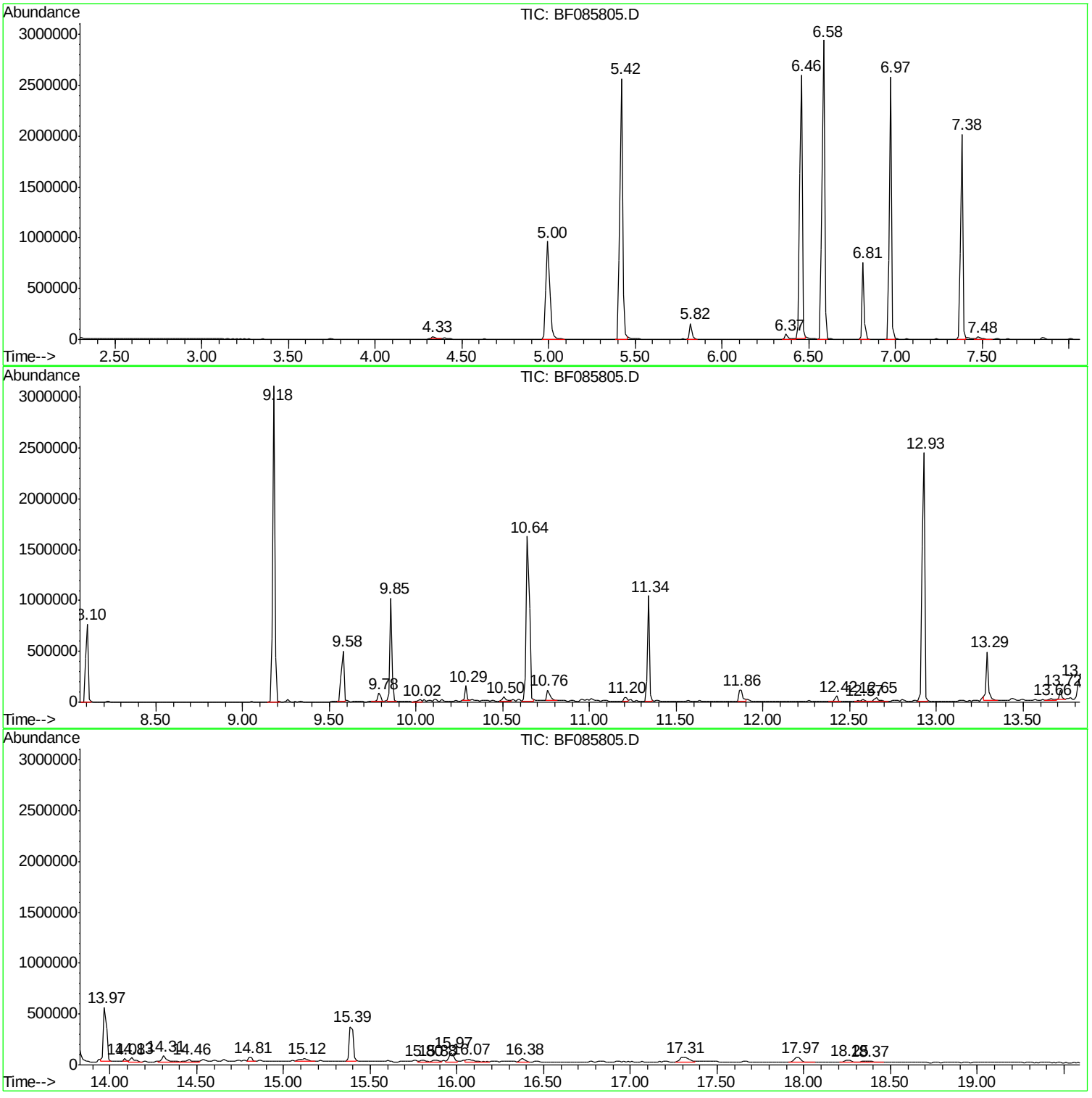
SP-8

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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Operator : UM/SJ
Sample : H1941-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

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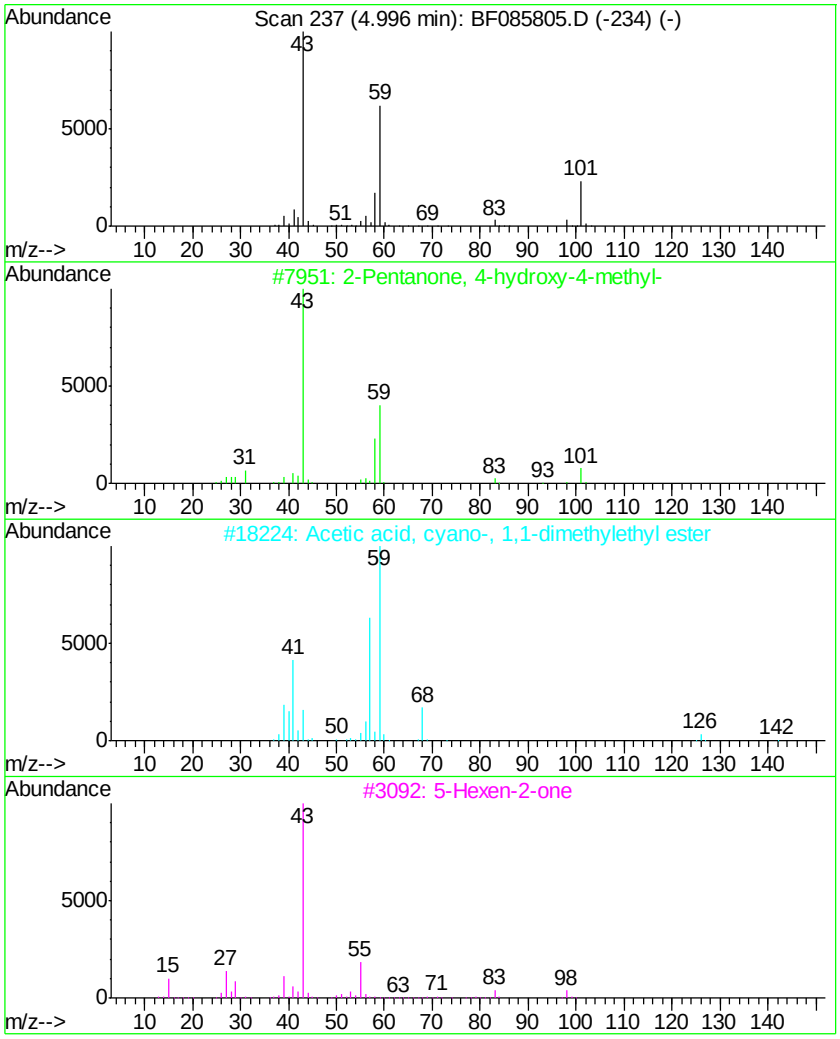
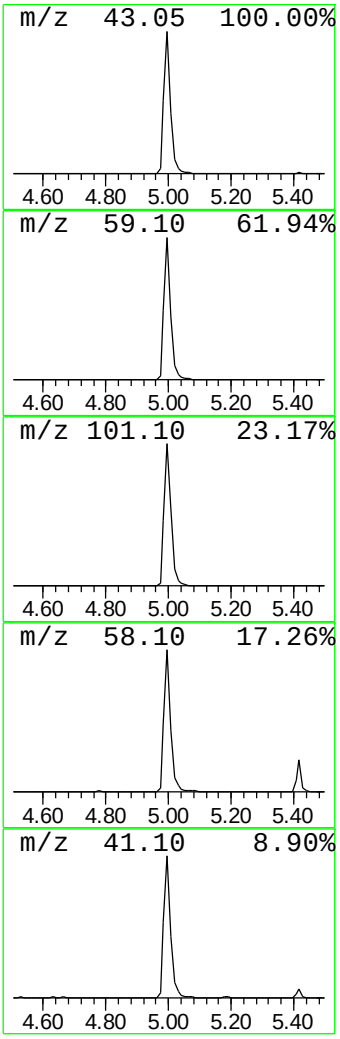
Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.00	43.50 ng	1415890	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, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	23
3	5-Hexen-2-one	98	C6H10O	000109-49-9	9
4	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9





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Instrument : BNA_F
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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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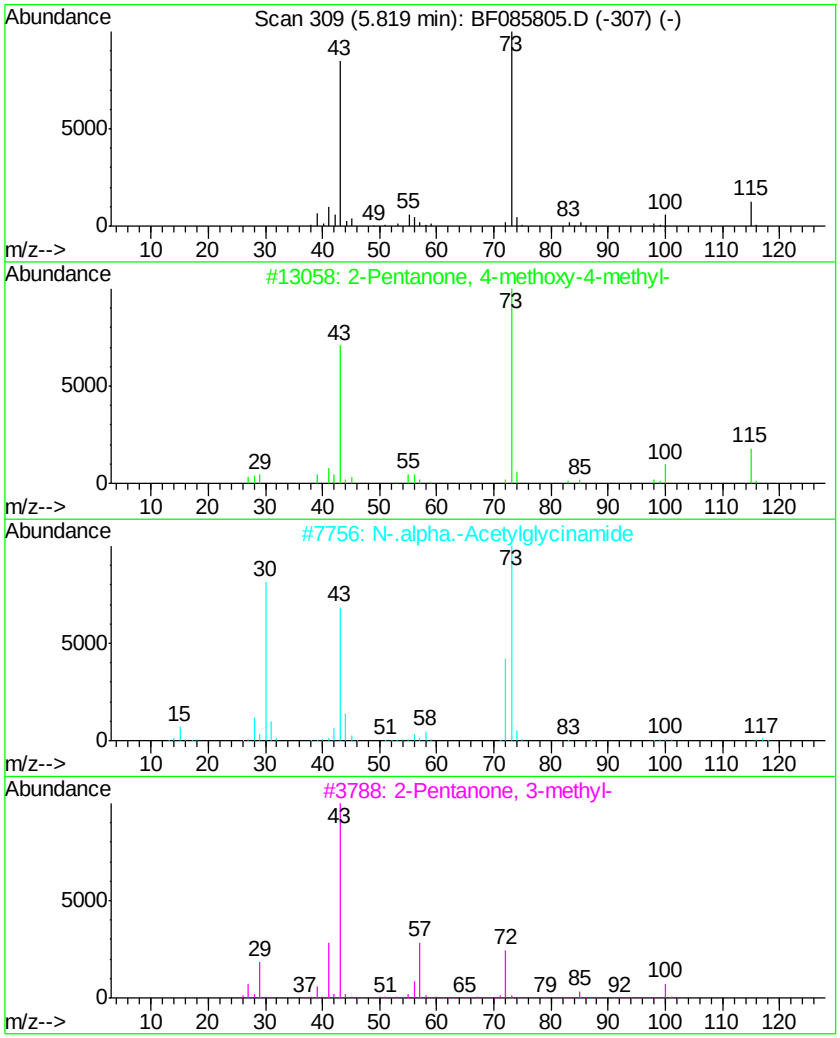
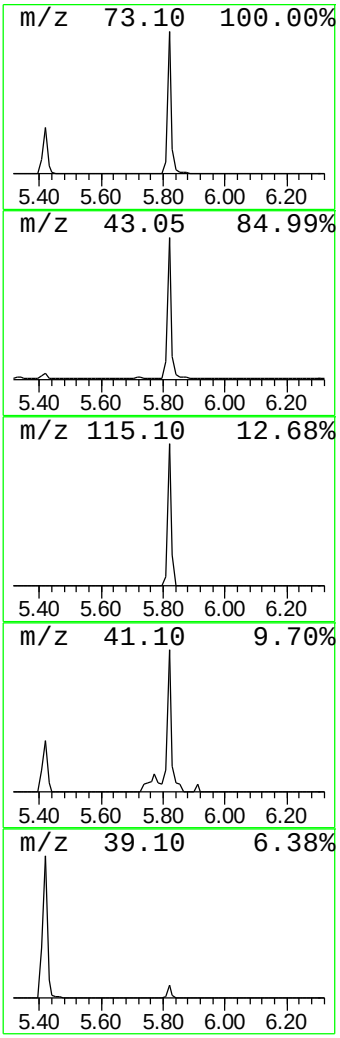
Peak Number 2 2-Pentanone, 4-methoxy-4-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.82	4.15 ng	134993	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	2-Pentanone, 3-methyl-	100	C6H12O	000565-61-7	27
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5	1,3-Dioxolane, 2-(1-methylethyl)-	116	C6H12O2	000822-83-3	10





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Data File : BF085805.D
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Operator : UM/SJ
Sample : H1941-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-8

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

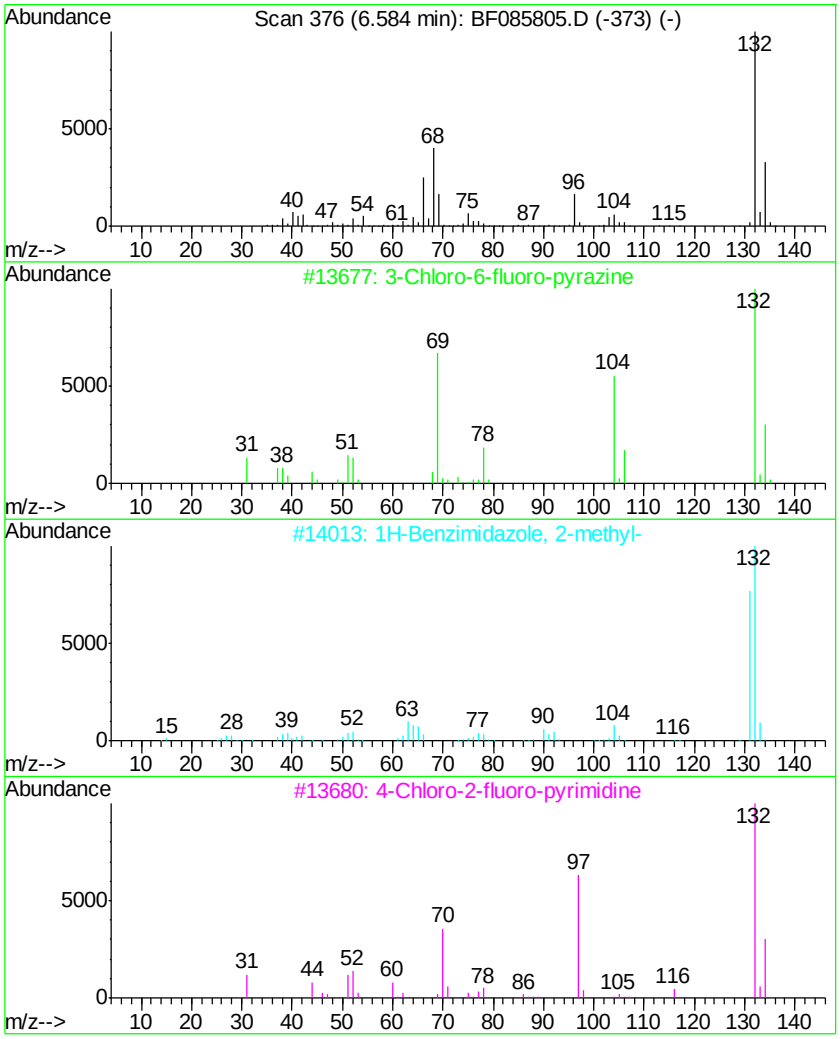
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

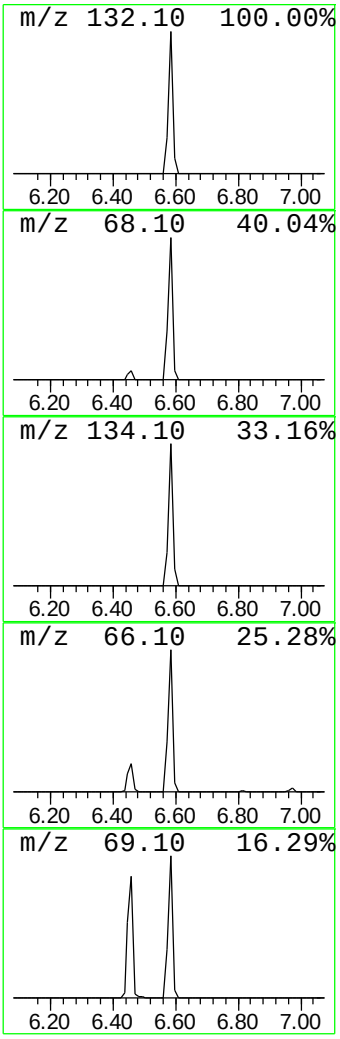
****

Peak Number 3 unknown6.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	86.43 ng	2813170	1,4-Dichlorobenzene-d4	6.81

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	10
3	4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
4	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	9
5	1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9





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Sample : H1941-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-8

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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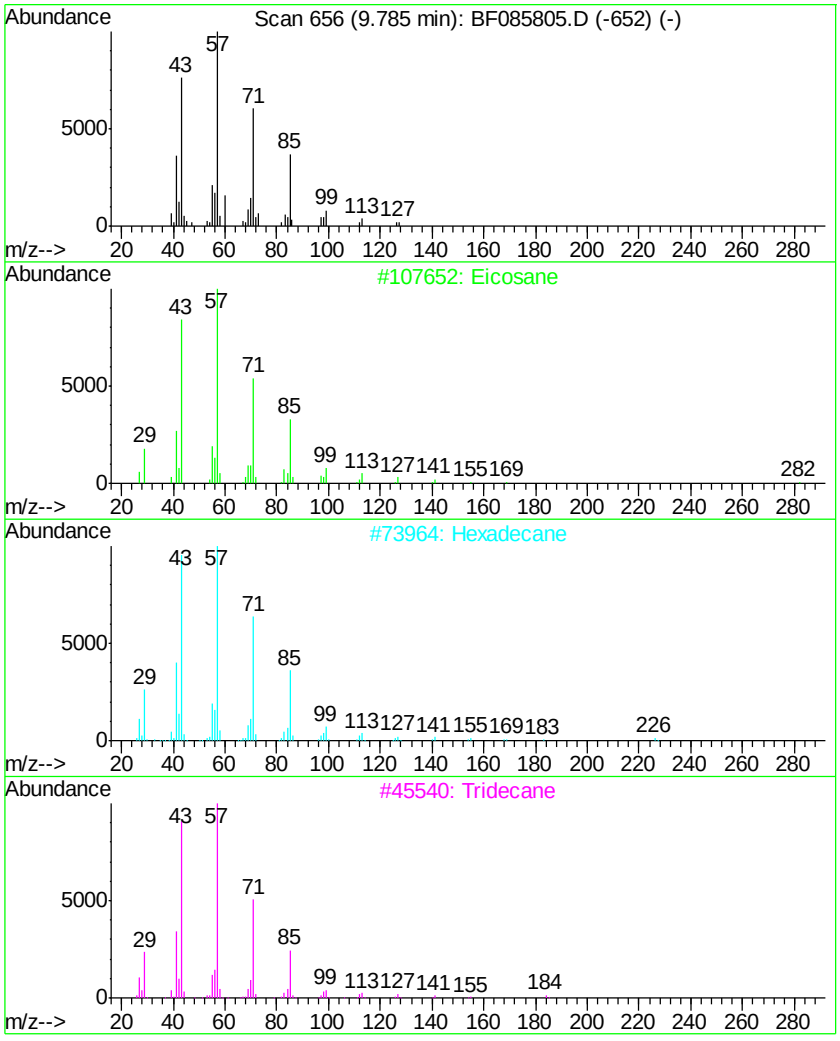
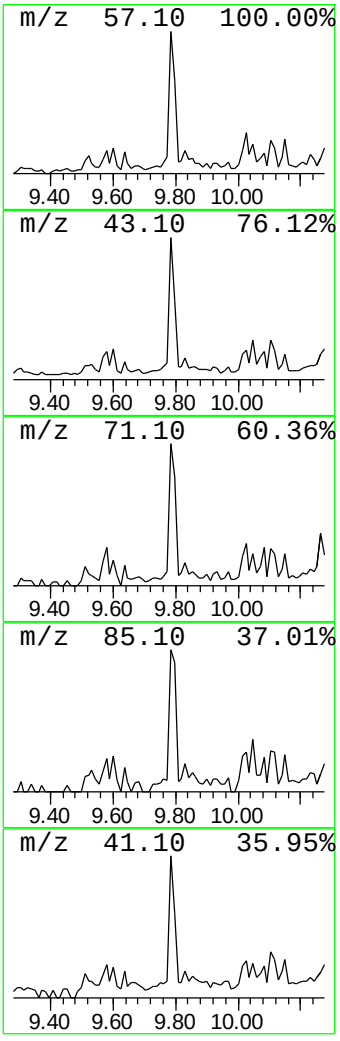
Peak Number 5 Eicosane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.78	2.52 ng	115064	Acenaphthene-d10	9.85

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	72
2	Hexadecane	226	C16H34	000544-76-3	72
3	Tridecane	184	C13H28	000629-50-5	64
4	Tetradecane	198	C14H30	000629-59-4	64
5	Pentadecane	212	C15H32	000629-62-9	59





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Sample : H1941-02
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ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF032416.M
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TIC Library : C:\DATABASE\NIST02.L
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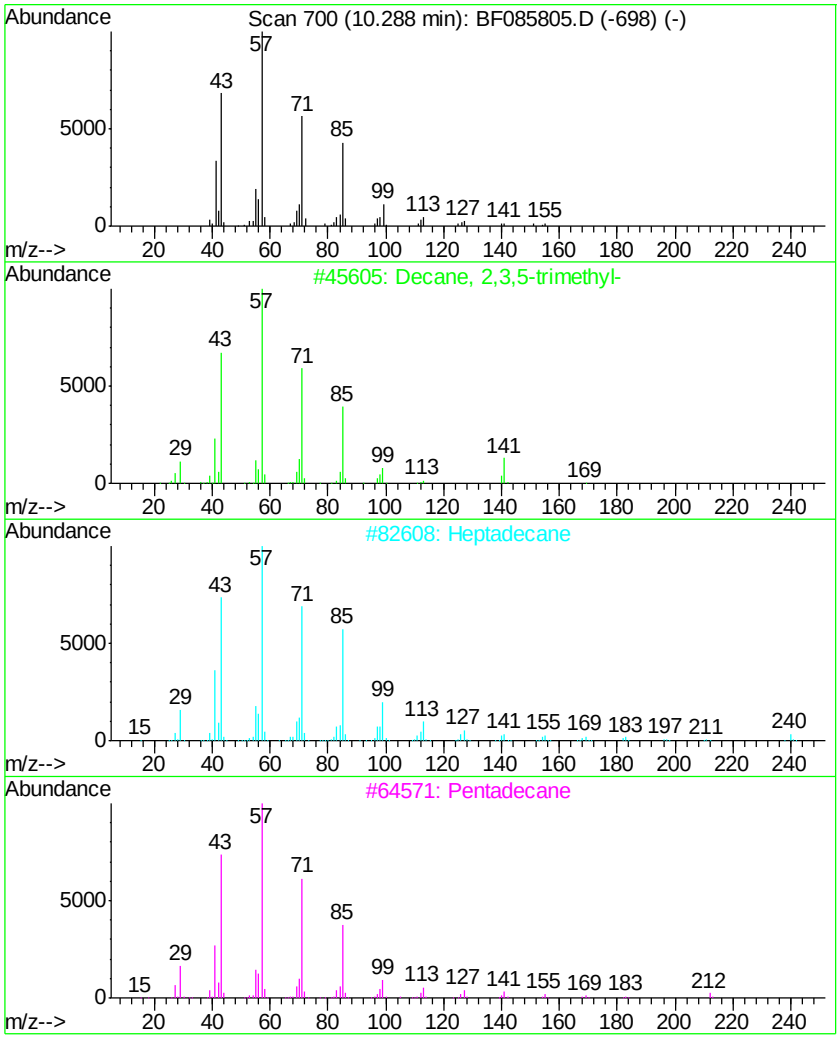
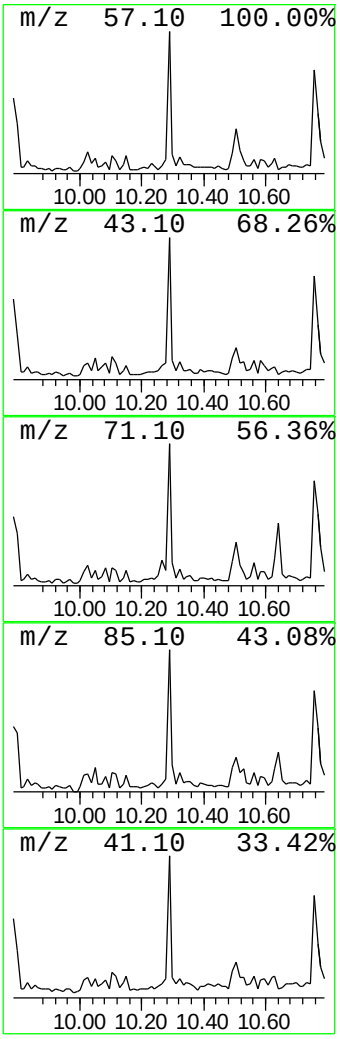
Peak Number 6 Decane, 2,3,5-trimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.29	2.55 ng	116465	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	Heptadecane	240	C17H36	000629-78-7	90
3	Pentadecane	212	C15H32	000629-62-9	78
4	Tetradecane	198	C14H30	000629-59-4	72
5	Decane, 3-bromo-	220	C10H21Br	030571-71-2	72





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ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-8

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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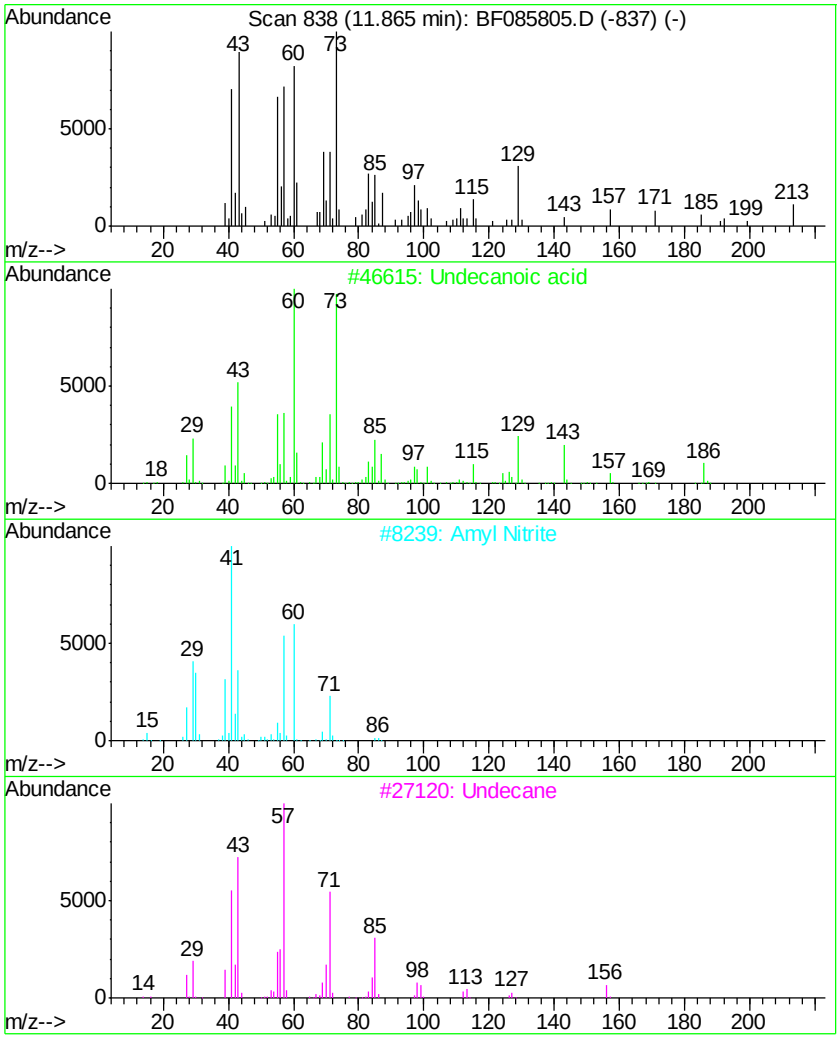
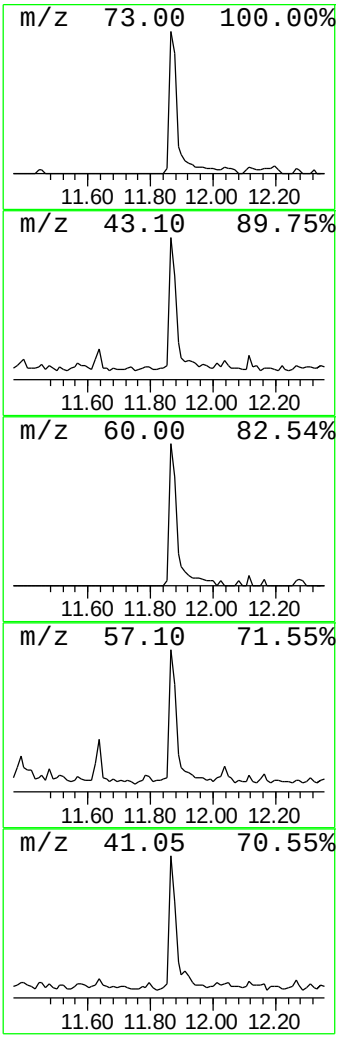
Peak Number 8 Undecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.86	3.99 ng	178021	Phenanthrene-d10	11.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecanoic acid	186	C11H22O2	000112-37-8	43
2	Amyl Nitrite	117	C5H11NO2	000110-46-3	38
3	Undecane	156	C11H24	001120-21-4	11
4	1-Tetradecanol, acetate	256	C16H32O2	000638-59-5	11
5	Heptadecane, 2-methyl-	254	C18H38	001560-89-0	10





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

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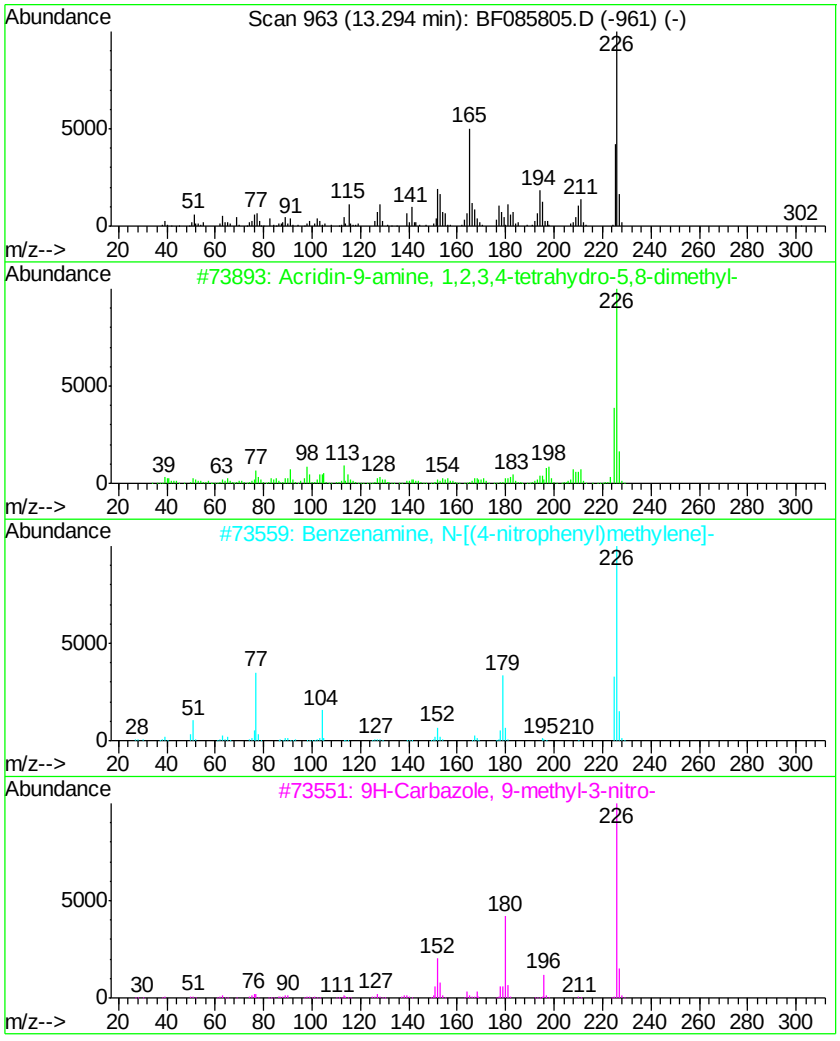
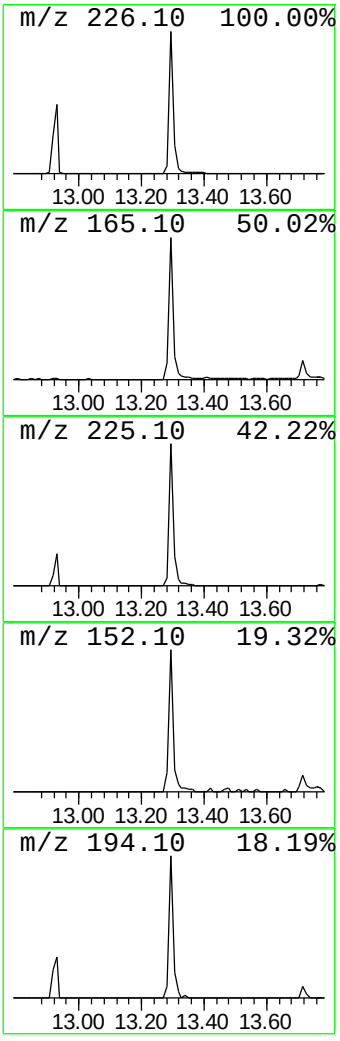
Peak Number 9 Acridin-9-amine, 1,2,3,4-te... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.29	14.61 ng	449228	Chrysene-d12	13.97

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acridin-9-amine, 1,2,3,4-tetrahy...	226	C15H18N2	297758-19-1	64
2	Benzenamine, N-[(4-nitrophenyl)m...	226	C13H10N2O2	000785-80-8	46
3	9H-Carbazole, 9-methyl-3-nitro-	226	C13H10N2O2	061166-05-0	45
4	1-Ethyl-4-methoxy-9H-pyrido[3,4-...	226	C14H14N2O	026585-14-8	43
5	2-Furfurylidene-7-hydroxy-1-inda...	226	C14H10O3	1000244-80-4	41





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Operator : UM/SJ
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Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-8

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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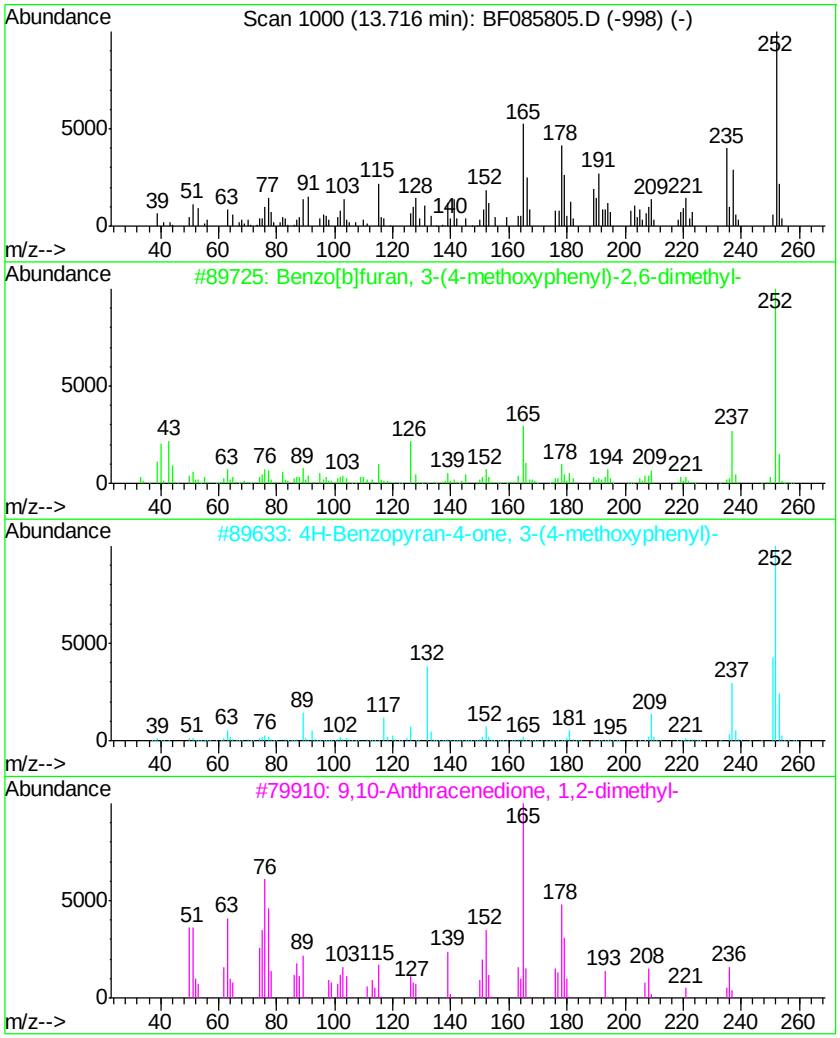
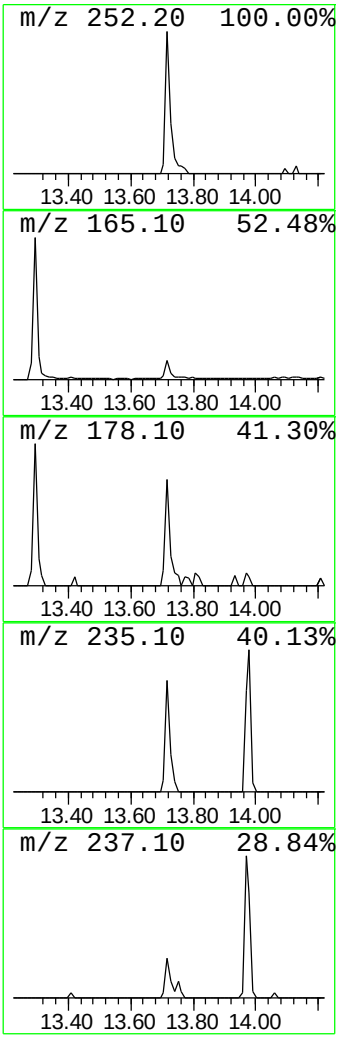
Peak Number 10 unknown13.72 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.72	3.16 ng	97114	Chrysene-d12	13.97

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]furan, 3-(4-methoxypheny...	252	C17H16O2	074228-98-1	47
2	4H-Benzopyran-4-one, 3-(4-methox...	252	C16H12O3	001621-58-5	46
3	9,10-Anthracenedione, 1,2-dimethyl-	236	C16H12O2	003285-98-1	44
4	2-Propenoic acid, 3-(4-phenylphe...	252	C17H16O2	040795-99-1	43
5	Adamantane-2,6-dione, bis(ethyle...	252	C14H20O4	060797-89-9	42





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Sample : H1941-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-8

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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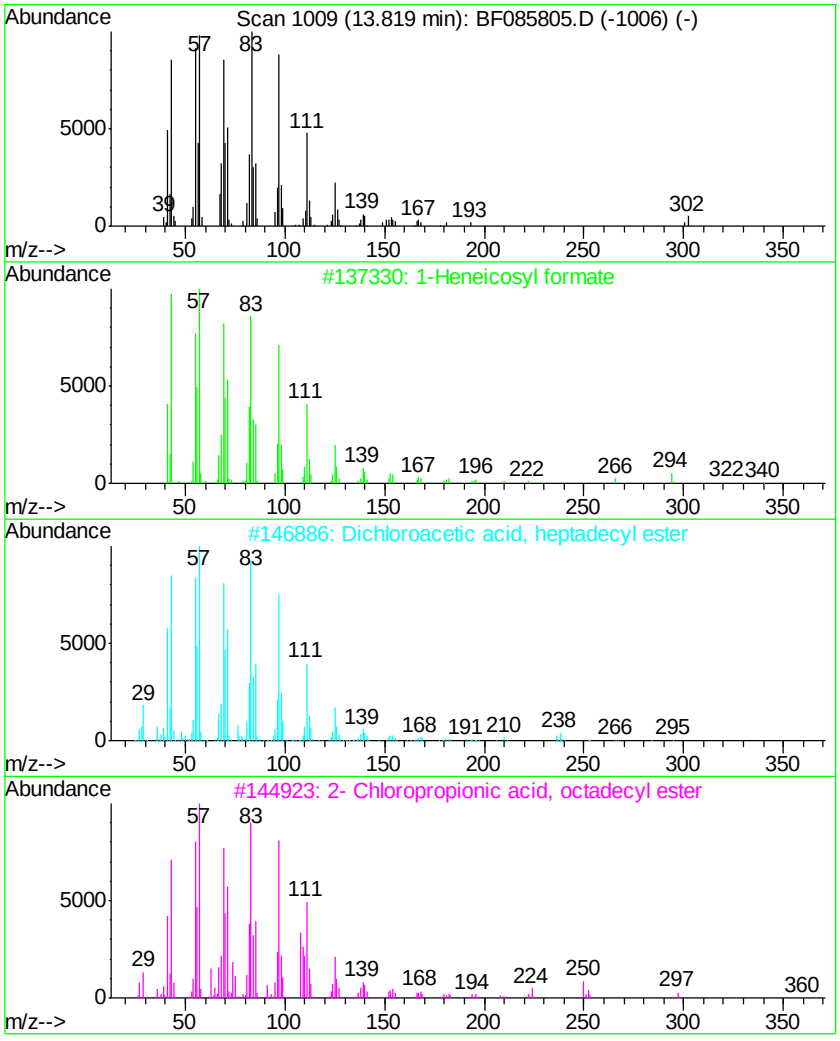
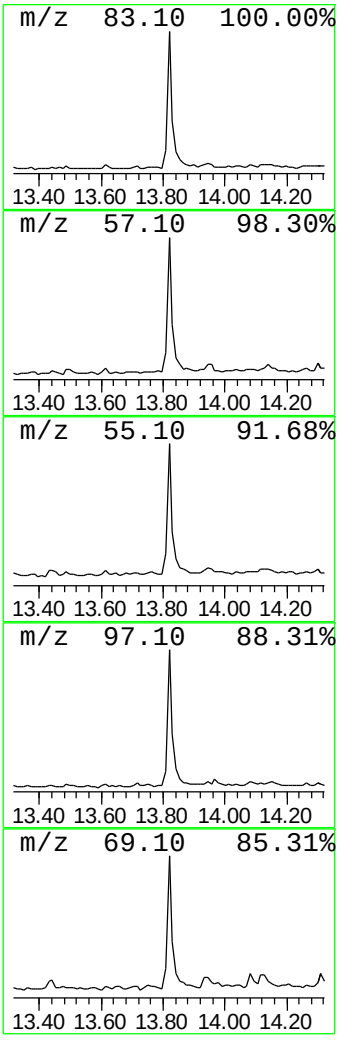
Peak Number 11 1-Heneicosyl formate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.82	8.18 ng	251719	Chrysene-d12	13.97

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosyl formate	340	C22H44O2	077899-03-7	95
2	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	95
3	2- Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	94
4	1-Nonadecanol	284	C19H40O	001454-84-8	93
5	1-Nonadecene	266	C19H38	018435-45-5	91





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Acq On : 28 Mar 2016 16:57
Operator : UM/SJ
Sample : H1941-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-8

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
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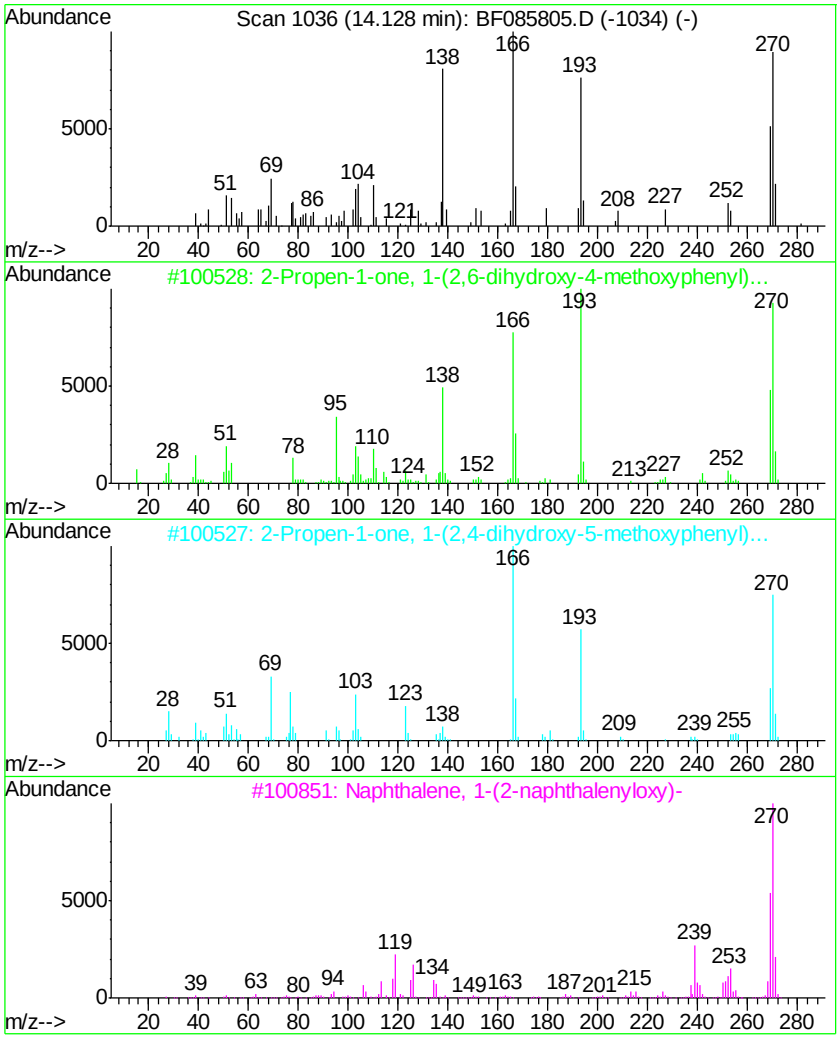
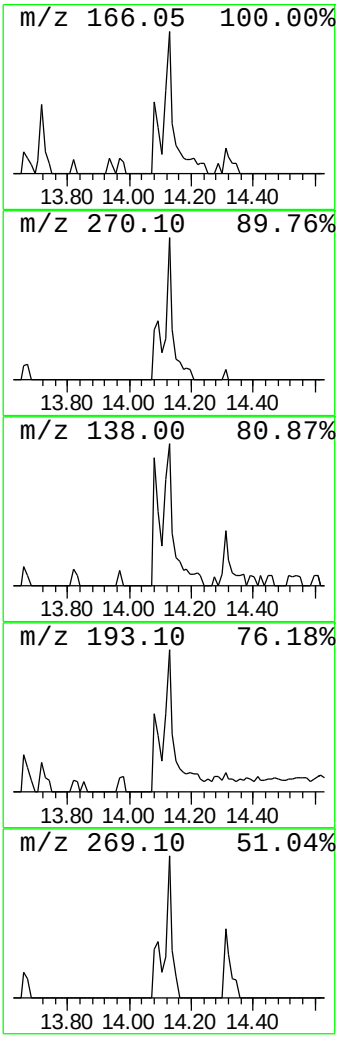
Peak Number 12 2-Propen-1-one, 1-(2,6-dihy... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.13	2.44 ng	75070	Chrysene-d12	13.97

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propen-1-one, 1-(2,6-dihydroxy...	270	C16H14O4	018956-15-5	58
2	2-Propen-1-one, 1-(2,4-dihydroxy...	270	C16H14O4	028143-82-0	53
3	Naphthalene, 1-(2-naphthalenyloxy)-	270	C20H14O	000611-49-4	22
4	Benzene, 1,2-methylenedioxy-4-(3...	270	C14H10N2O4	191595-07-0	22
5	4-(6-Isopropyl-4-methyl-1-naphth...	270	C18H22O2	094265-06-2	14





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ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-8

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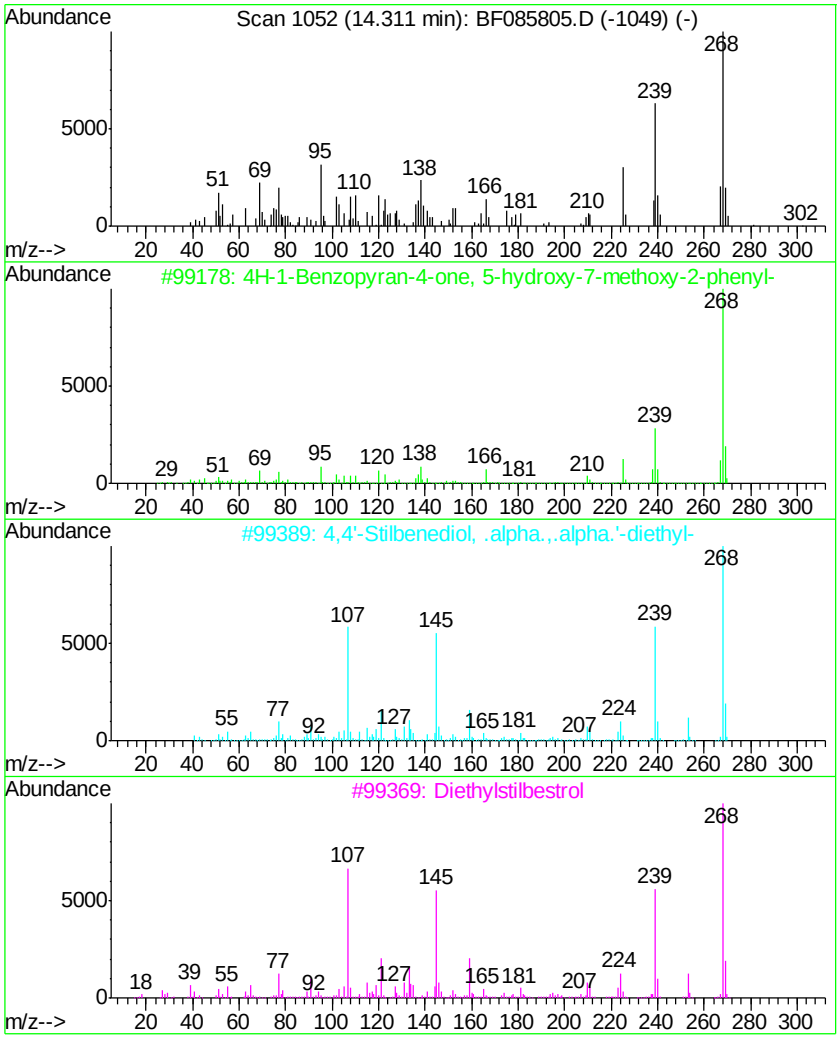
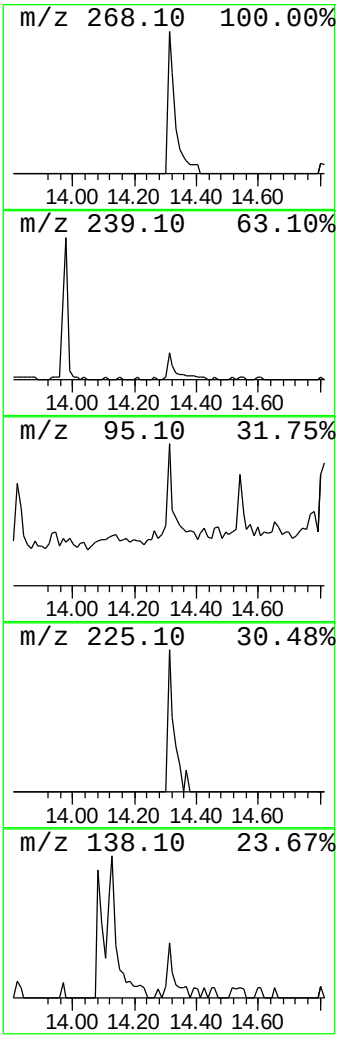
Peak Number 13 4H-1-Benzopyran-4-one, 5-hy... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.31	3.02 ng	92932	Chrysene-d12	13.97

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-1-Benzopyran-4-one, 5-hydroxy...	268	C16H12O4	000520-28-5	91
2	4,4'-Stilbenediol, .alpha.,.alph...	268	C18H20O2	006898-97-1	52
3	Diethylstilbestrol	268	C18H20O2	000056-53-1	52
4	Dinaphtho[2,1-b:1',2'-d]furan	268	C20H12O	000194-63-8	50
5	N4,N4'-Dipropyl-biphenyl-4,4'-di...	268	C18H24N2	017576-22-6	50





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Misc :
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Instrument : BNA_F
ClientSampleId : SP-8

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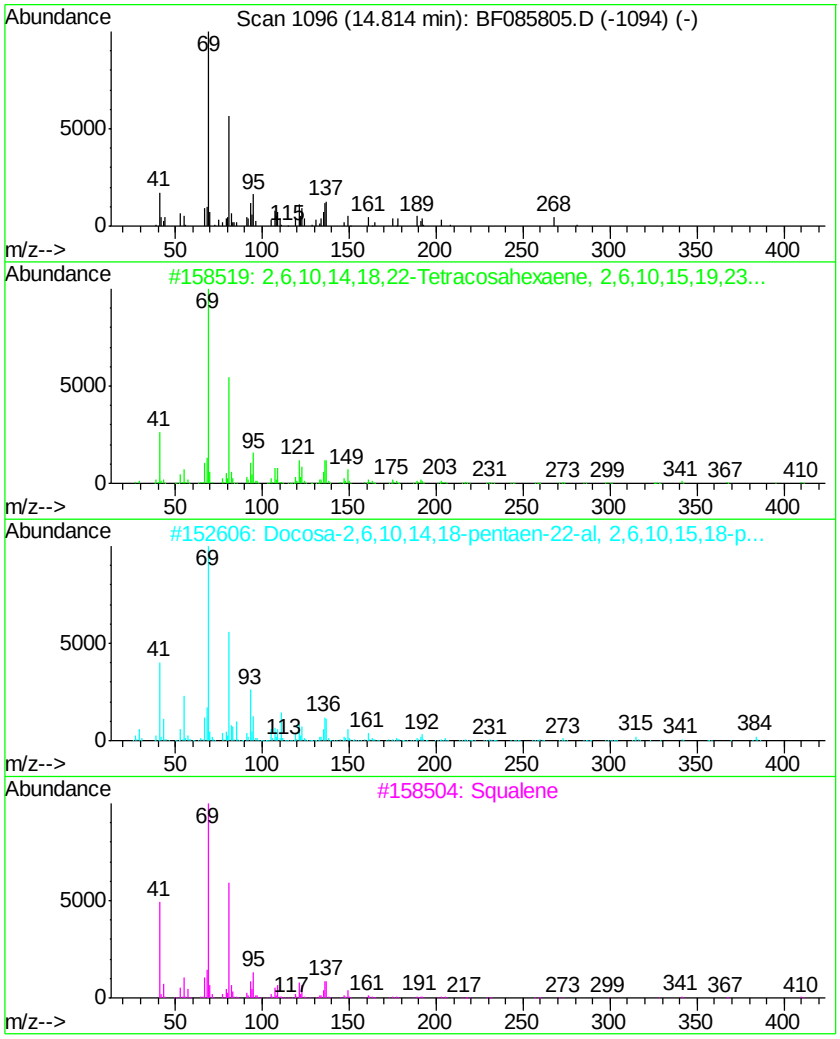
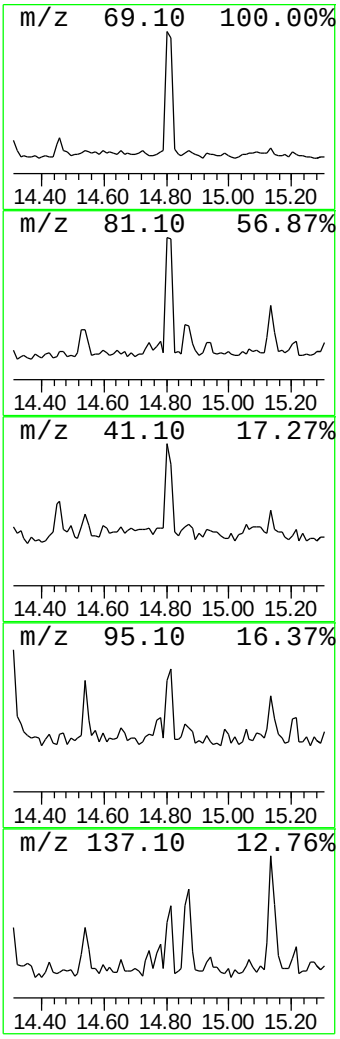
Peak Number 14 2,6,10,14,18,22-Tetracosahe... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.81	2.24 ng	56685	Perylene-d12	15.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6,10,14,18,22-Tetracosahexaene...	410	C30H50	000111-02-4	87
2	Docosa-2,6,10,14,18-pentaen-22-a...	384	C27H44O	1000163-04-7	86
3	Squalene	410	C30H50	007683-64-9	83
4	4,9,13,17-Tetramethyl-4,8,12,16-...	316	C22H36O	056882-09-8	80
5	Hexadeca-2,6,10,14-tetraen-1-ol,...	290	C20H34O	007614-21-3	72





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Misc :
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ClientSampleId : SP-8

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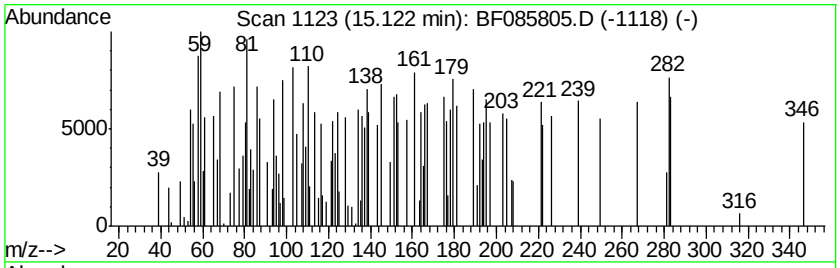
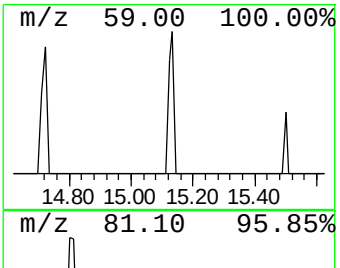
Peak Number 15 unknown15.12 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.12	3.99 ng	101079	Perylene-d12	15.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyridine, 2,6-dichloro-3-nitro-	192	C5H2Cl2N2O2	016013-85-7	25
2	Carbazol-2-yl methyl ketone thio...	282	C15H14N4S	1000254-48-4	22
3	Methanone, bis(5-methyl-1H-1,2,3...	192	C7H8N6O	051719-73-4	11
4	Phenanthrene, 9,10-dihydro-1-met...	194	C15H14	095676-48-5	11
5	1-Oxaspiro[2.5]octane, 4,4-dimet...	194	C13H22O	054345-65-2	10





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ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-8

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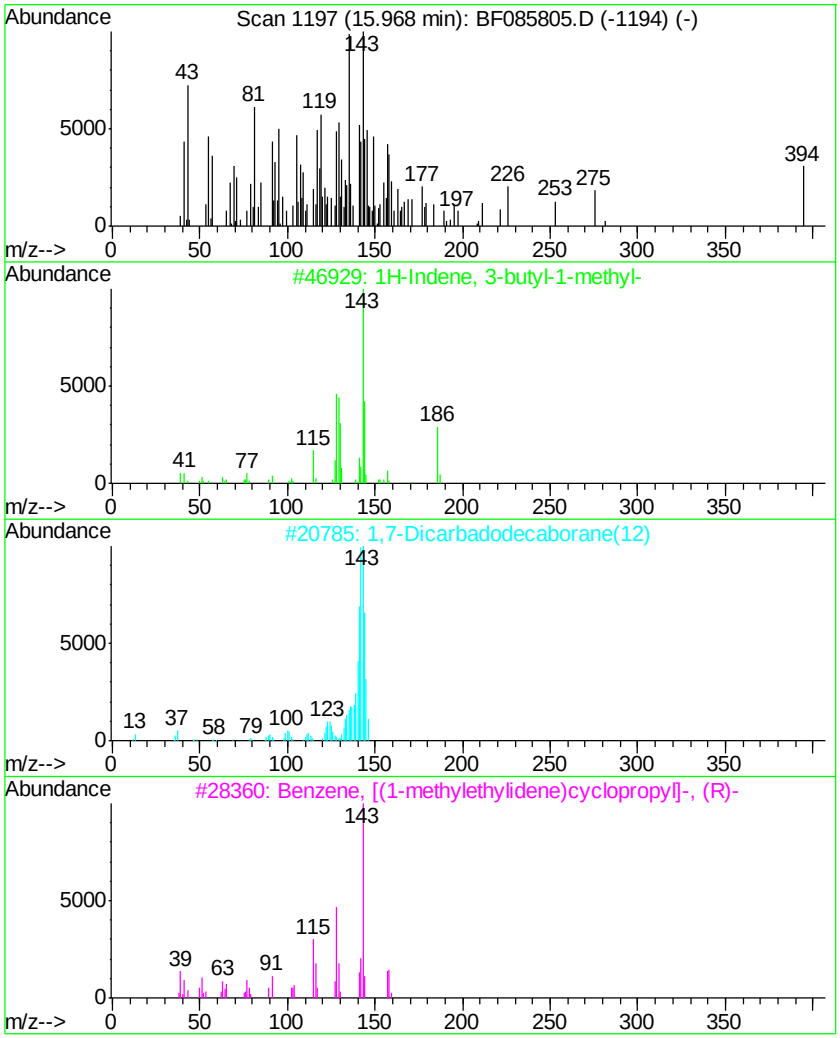
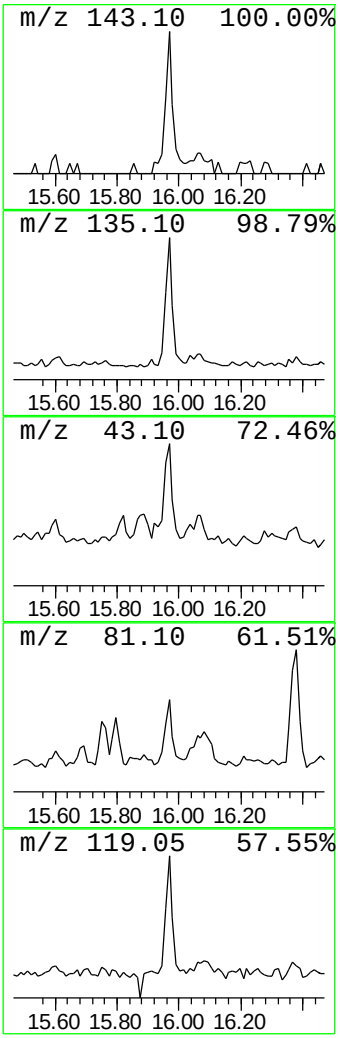
Peak Number 16 unknown15.97 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.97	6.22 ng	157463	Perylene-d12	15.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 3-butyl-1-methyl-	186	C14H18	111400-84-1	27
2	1,7-Dicarbadodecaborane(12)	146	C2H12B10	016986-24-6	25
3	Benzene, [(1-methylethylidene)cy...	158	C12H14	024524-55-8	22
4	Quinoline, 5-methyl-	143	C10H9N	007661-55-4	18
5	Quinoline, 6-methyl-	143	C10H9N	000091-62-3	18





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Sample : H1941-02
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SP-8

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TIC Library : C:\DATABASE\NIST02.L
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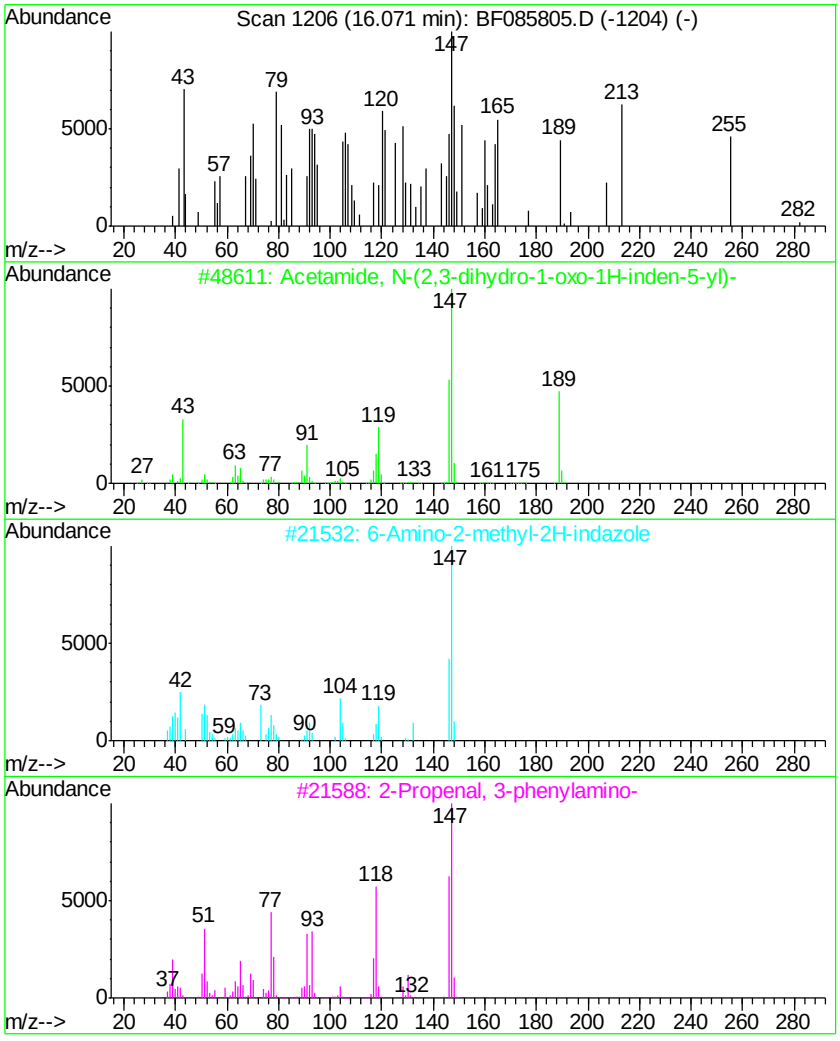
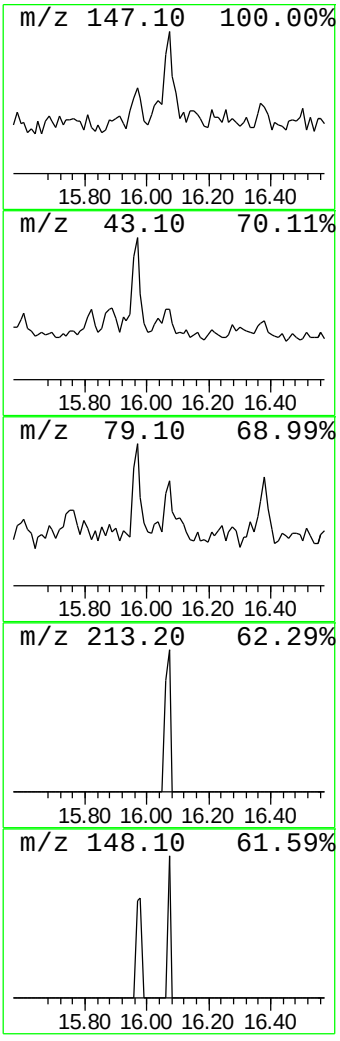
Peak Number 17 unknown16.07 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.07	2.88 ng	72905	Perylene-d12	15.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetamide, N-(2,3-dihydro-1-oxo-...	189	C11H11NO2	058161-35-6	22
2	6-Amino-2-methyl-2H-indazole	147	C8H9N3	050593-30-1	9
3	2-Propenal, 3-phenylamino-	147	C9H9NO	025299-39-2	9
4	Pyridine, 3,5-dichloro-	147	C5H3Cl2N	002457-47-8	9
5	1-(2'-Hydroxy-5'-methylphenyl)-1...	165	C9H11NO2	1000146-89-9	9





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Instrument : BNA_F
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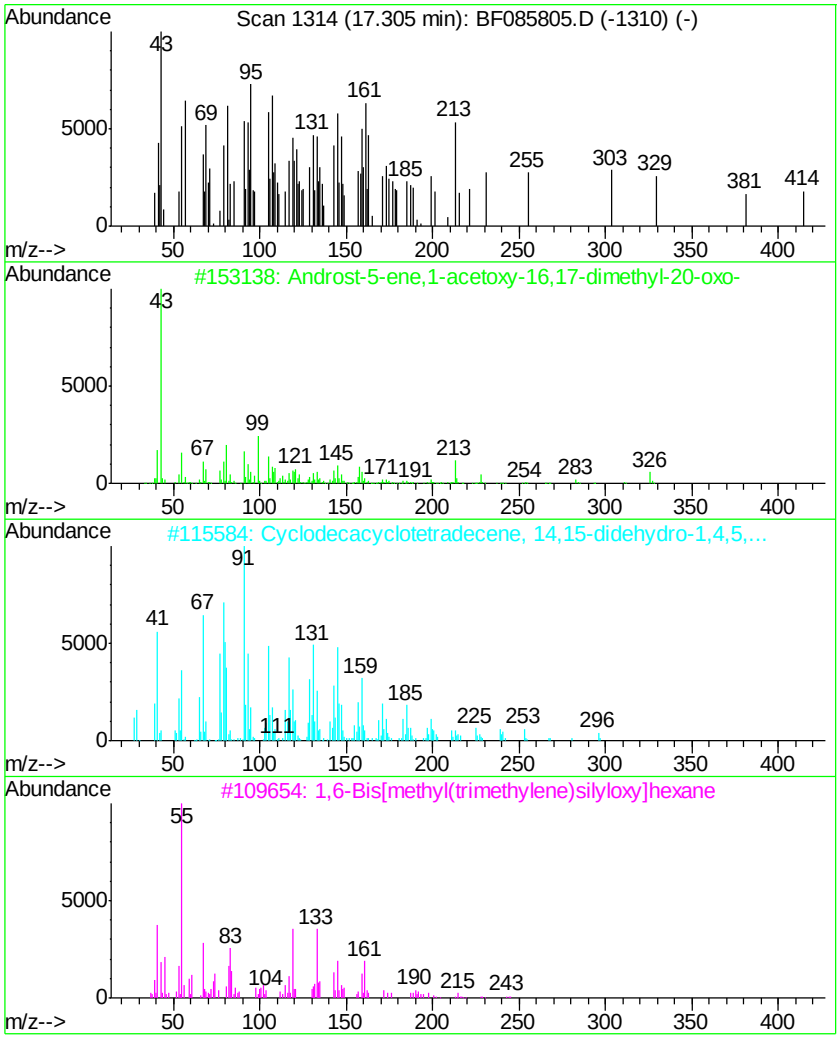
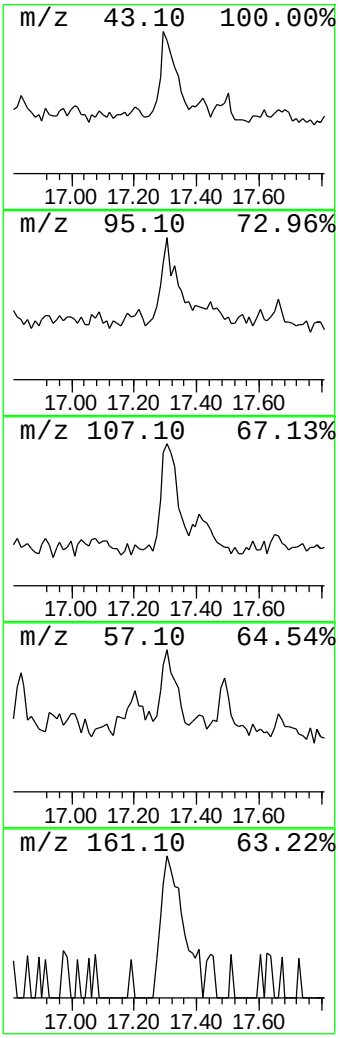
Peak Number 18 unknown17.31 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.31	6.39 ng	161815	Perylene-d12	15.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Androst-5-ene,1-acetoxy-16,17-di...	386	C25H38O3	294866-91-4	35
2	Cyclodecacyclotetradecene, 14,15...	296	C22H32	014113-61-2	22
3	1,6-Bis[methyl(trimethylene)sily...	286	C14H30O2Si2	1000217-00-5	17
4	3-(3,7-Dimethyl-octa-2,6-dienyl)...	258	C17H22O2	1000192-28-0	10
5	N-(3-Chloro-4-fluoro-phenyl)-2-(...	329	C14H10Cl2FNOS	1000295-12-7	10





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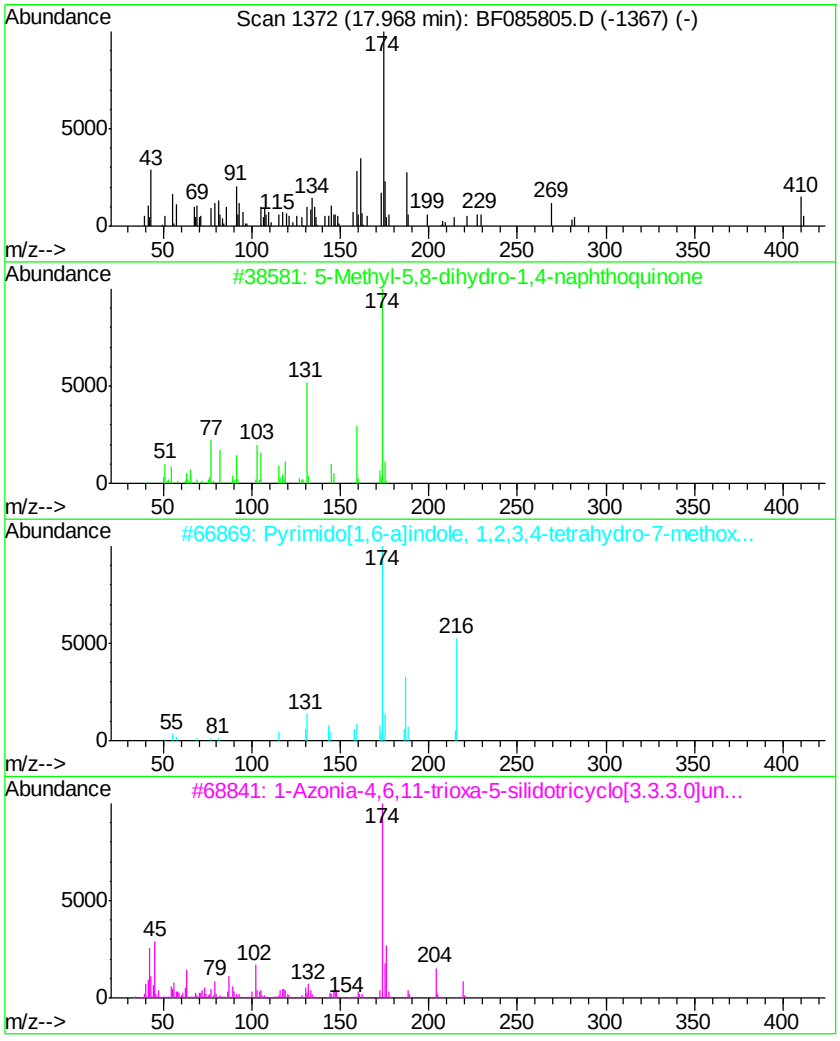
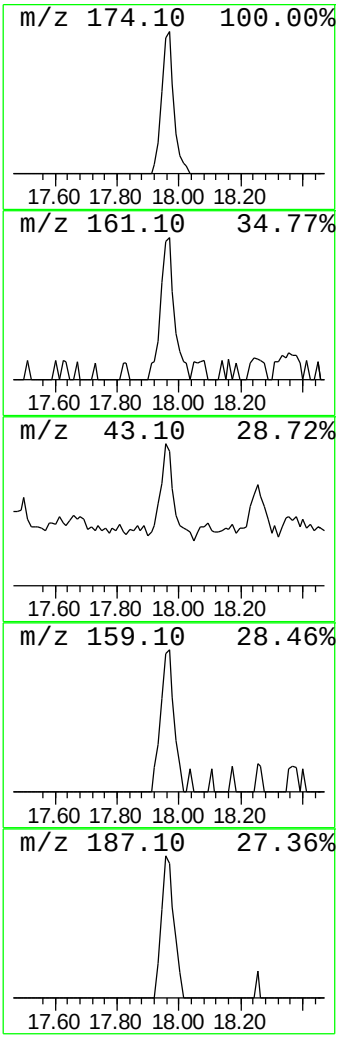
Peak Number 19 unknown17.97 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.97	5.03 ng	127324	Perylene-d12	15.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Methyl-5,8-dihydro-1,4-naphtho...	174	C11H10O2	086759-40-2	43
2	Pyrimido[1,6-a]indole, 1,2,3,4-t...	216	C13H16N2O	038349-09-6	43
3	1-Azonia-4,6,11-trioxa-5-silidot...	219	C8H17NO4Si	349498-09-5	38
4	6-Indolizinecarboxamide, 2-methyl-	174	C10H10N2O	022380-20-7	35
5	8-Amino-6-methoxyquinoline	174	C10H10N2O	000090-52-8	32





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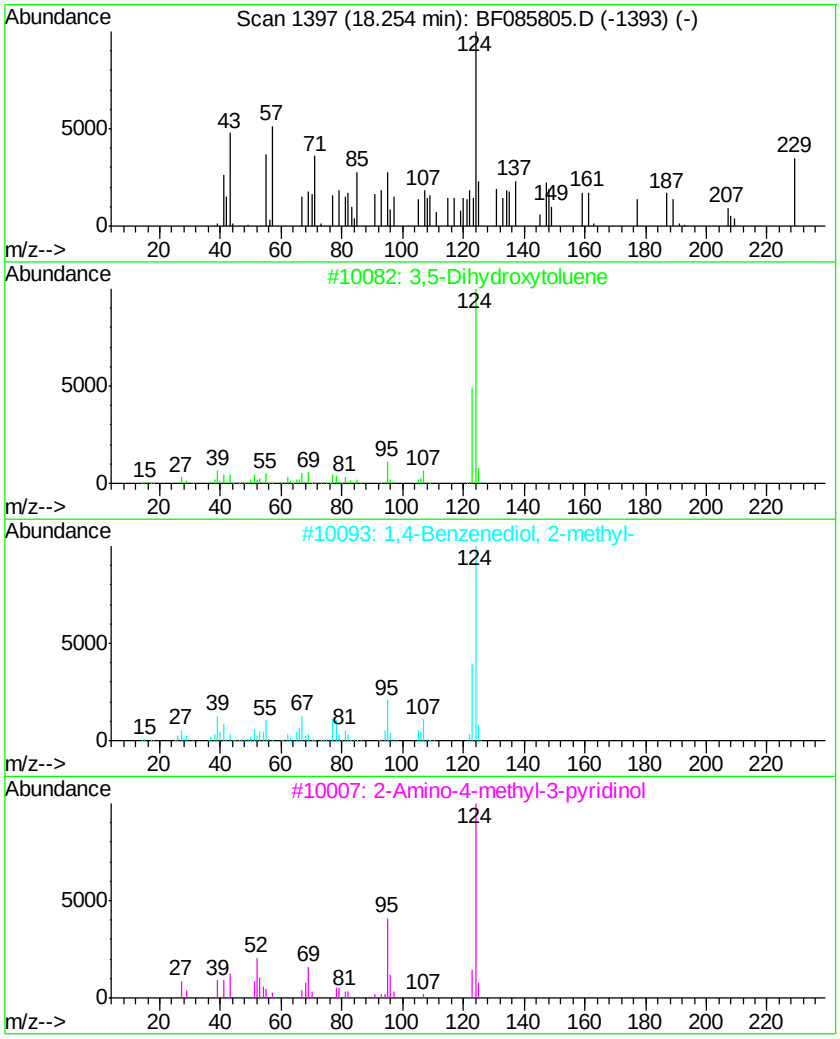
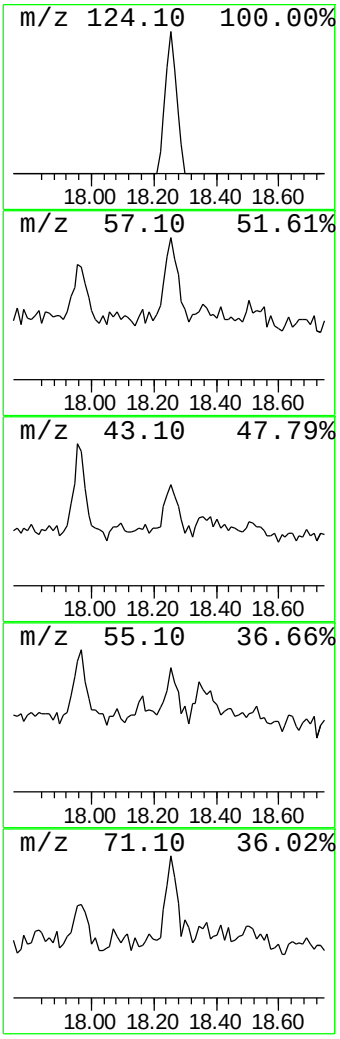
Peak Number 20 unknown18.25 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.25	2.21 ng	55894	Perylene-d12	15.40

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,5-Dihydroxytoluene	124	C7H8O2	000504-15-4	35
2	1,4-Benzenediol, 2-methyl-	124	C7H8O2	000095-71-6	35
3	2-Amino-4-methyl-3-pyridinol	124	C6H8N2O	020348-18-9	35
4	4,5-Dimethyl-2-pyrimidone	124	C6H8N2O	034939-17-8	27
5	Pyrazole, 3-methylaminomethyl-5-...	153	C8H15N3	1000262-50-0	27





Tentatively Identified Compound (LSC) summary

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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	43.5	ng	1415890	1	6.81	650991	20.0
2-Pentanone, 4-me...	5.82	4.2	ng	134993	1	6.81	650991	20.0
unknown6.58	6.58	86.4	ng	2813170	1	6.81	650991	20.0
Eicosane	9.78	2.5	ng	115064	3	9.85	914555	20.0
Decane, 2,3,5-tri...	10.29	2.5	ng	116465	3	9.85	914555	20.0
Undecanoic acid	11.86	4.0	ng	178021	4	11.34	892672	20.0
Acridin-9-amine, ...	13.29	14.6	ng	449228	5	13.97	615132	20.0
unknown13.72	13.72	3.2	ng	97114	5	13.97	615132	20.0
1-Heneicosyl formate	13.82	8.2	ng	251719	5	13.97	615132	20.0
2-Propen-1-one, 1...	14.13	2.4	ng	75070	5	13.97	615132	20.0
4H-1-Benzopyran-4...	14.31	3.0	ng	92932	5	13.97	615132	20.0
2,6,10,14,18,22-T...	14.81	2.2	ng	56685	6	15.40	506208	20.0
unknown15.12	15.12	4.0	ng	101079	6	15.40	506208	20.0
unknown15.97	15.97	6.2	ng	157463	6	15.40	506208	20.0
unknown16.07	16.07	2.9	ng	72905	6	15.40	506208	20.0
unknown17.31	17.31	6.4	ng	161815	6	15.40	506208	20.0
unknown17.97	17.97	5.0	ng	127324	6	15.40	506208	20.0
unknown18.25	18.25	2.2	ng	55894	6	15.40	506208	20.0