LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
Data File : BF099186.D
Acq On : 7 Oct 2017 11:26
Operator : SJ/JU
Sample : I5586-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G72-0-1

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-BF092317.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.158	8	12	29	rVB	164662	229570	5.90%	0.635%
2	3.463	232	234	251	rBV2	12977	46798	1.20%	0.130%
3	5.093	507	511	523	rVB	434765	635309	16.33%	1.758%
4	5.493	574	579	582	rBV	2996725	3180565	81.76%	8.801%
5	6.504	740	751	754	rBV	2849337	3391703	87.19%	9.386%
6	6.640	769	774	777	rBV	3269119	3313773	85.19%	9.170%
7	6.863	808	812	816	rBV	976464	786297	20.21%	2.176%
8	6.922	818	822	827	rBV	53251	51347	1.32%	0.142%
9	7.022	834	839	842	rBV	2944469	2674760	68.76%	7.402%
10	7.428	902	908	911	rBV	1929955	2200067	56.56%	6.088%
11	7.516	917	923	928	rVB	35239	47530	1.22%	0.132%
12	8.145	1026	1030	1033	rBV	1280709	1066422	27.41%	2.951%
13	9.222	1207	1213	1216	rBV	3974292	3890072	100.00%	10.765%
14	9.486	1253	1258	1261	rBV2	72291	56943	1.46%	0.158%
15	9.610	1275	1279	1282	rVB	630664	484813	12.46%	1.342%
16	9.898	1321	1328	1332	rBV	1345765	1168772	30.04%	3.234%
17	10.098	1356	1362	1366	rBV4	40261	67055	1.72%	0.186%
18	10.322	1397	1400	1402	rVB2	62843	59006	1.52%	0.163%
19	10.533	1429	1436	1439	rBV3	31560	54241	1.39%	0.150%
20	10.692	1455	1463	1466	rBV	2013738	2036929	52.36%	5.637%
21	10.792	1477	1480	1486	rBV	81737	91325	2.35%	0.253%
22	11.039	1517	1522	1527	rVB3	32494	47826	1.23%	0.132%
23	11.239	1552	1556	1558	rBV2	63386	53082	1.36%	0.147%
24	11.269	1558	1561	1567	rVB3	21954	40082	1.03%	0.111%
25	11.386	1577	1581	1583	rBV	1238109	1065655	27.39%	2.949%
26	11.410	1583	1585	1588	rVV	294257	232644	5.98%	0.644%
27	11.457	1588	1593	1596	rVB2	54071	67237	1.73%	0.186%
28	11.592	1613	1616	1618	rBV	92256	88297	2.27%	0.244%
29	11.663	1626	1628	1634	rVB2	35695	40198	1.03%	0.111%
30	11.904	1661	1669	1674	rBV3	369582	504636	12.97%	1.396%
31	11.998	1682	1685	1687	rBV2	41764	45330	1.17%	0.125%
32	12.210	1718	1721	1726	rVB6	29085	44622	1.15%	0.123%
33	12.522	1771	1774	1778	rBV	47404	44527	1.14%	0.123%
34	12.598	1783	1787	1792	rBV	345483	322956	8.30%	0.894%

LSC Area Percent Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
Data File : BF099186.D
Acq On : 7 Oct 2017 11:26
Operator : SJ/JU
Sample : I5586-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G72-0-1

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-BF092317.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	12.686	1799	1802	1809	rBV2	92758	93072	2.39%	0.258%
36	12.827	1820	1826	1828	rBV	260341	269075	6.92%	0.745%
37	12.851	1828	1830	1841	rVB	393032	439438	11.30%	1.216%
38	12.969	1845	1850	1854	rVV	2720061	2895738	74.44%	8.013%
39	13.157	1879	1882	1885	rVB3	35535	45245	1.16%	0.125%
40	13.439	1925	1930	1938	rVB	648481	691695	17.78%	1.914%
41	13.851	1997	2000	2007	rVB2	304070	319176	8.20%	0.883%
42	14.022	2025	2029	2032	rBV	756210	775267	19.93%	2.145%
43	14.051	2032	2034	2040	rVB	108319	107549	2.76%	0.298%
44	14.480	2104	2107	2113	rBV3	94050	143425	3.69%	0.397%
45	14.645	2129	2135	2147	rBV4	193258	415002	10.67%	1.148%
46	15.063	2203	2206	2210	rBV	99654	152733	3.93%	0.423%
47	15.368	2255	2258	2264	rBV	58899	74750	1.92%	0.207%
48	15.433	2266	2269	2273	rVB	77796	95223	2.45%	0.264%
49	15.498	2275	2280	2284	rBV	556019	670015	17.22%	1.854%
50	15.651	2302	2306	2315	rVB2	116689	236440	6.08%	0.654%
51	16.974	2527	2531	2537	rBV2	37718	77054	1.98%	0.213%
52	17.204	2564	2570	2579	rVB4	115349	288163	7.41%	0.797%
53	17.551	2624	2629	2639	rVB8	74097	217569	5.59%	0.602%

Sum of corrected areas: 36137018

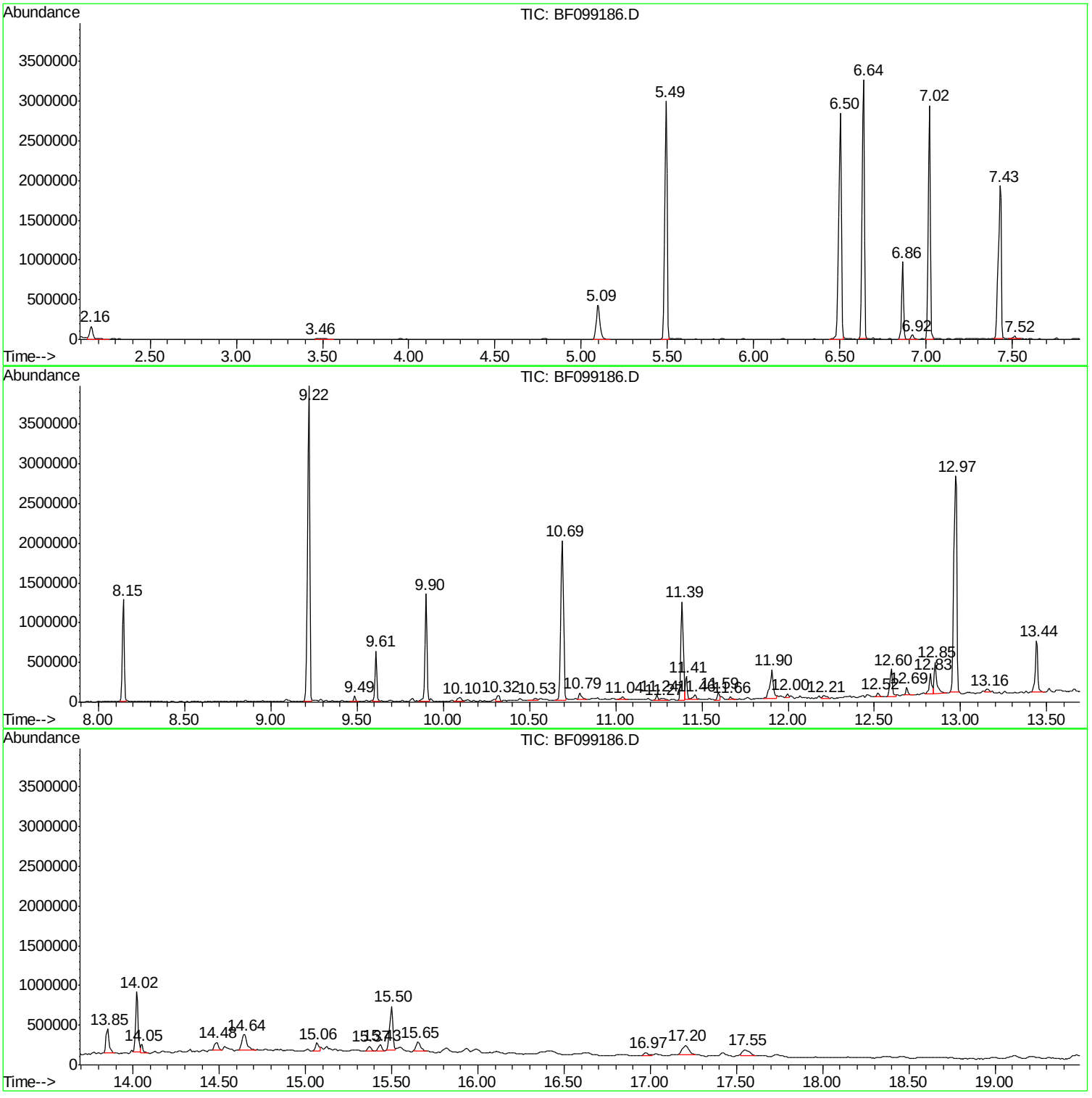
LSC Report - Integrated Chromatogram

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
Data File : BF099186.D
Acq On : 7 Oct 2017 11:26
Operator : SJ/JU
Sample : I5586-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G72-0-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
Data File : BF099186.D
Acq On : 7 Oct 2017 11:26
Operator : SJ/JU
Sample : I5586-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G72-0-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

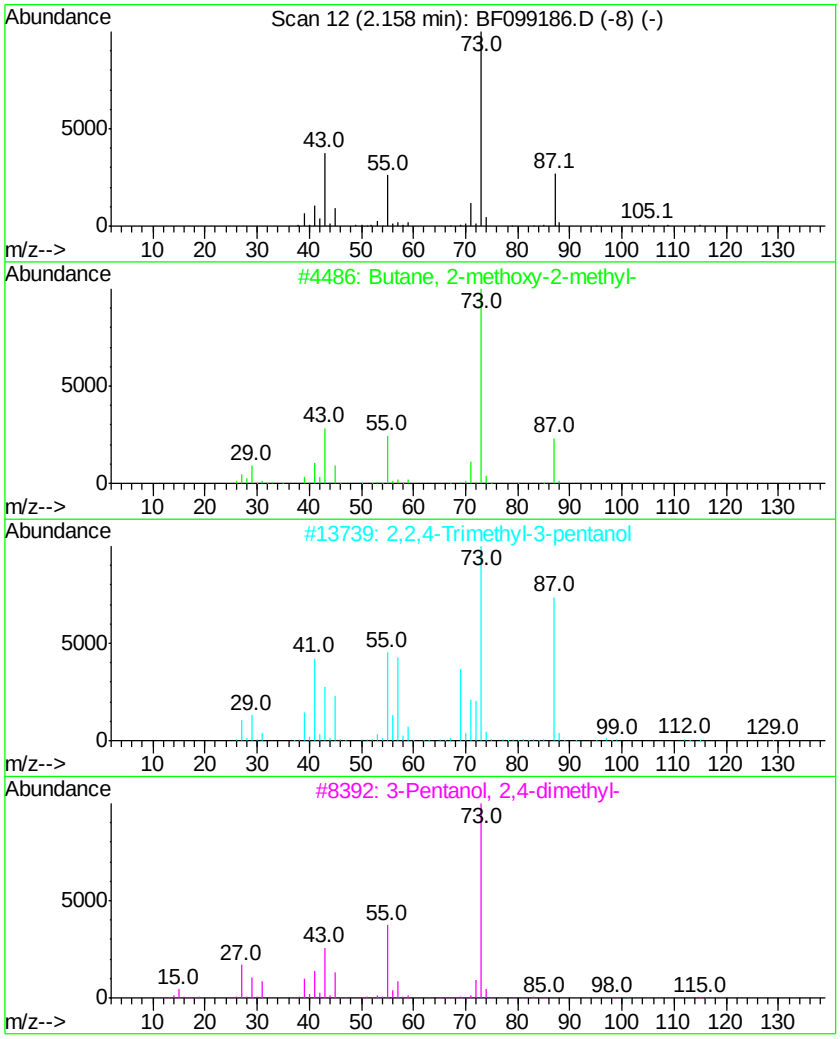
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

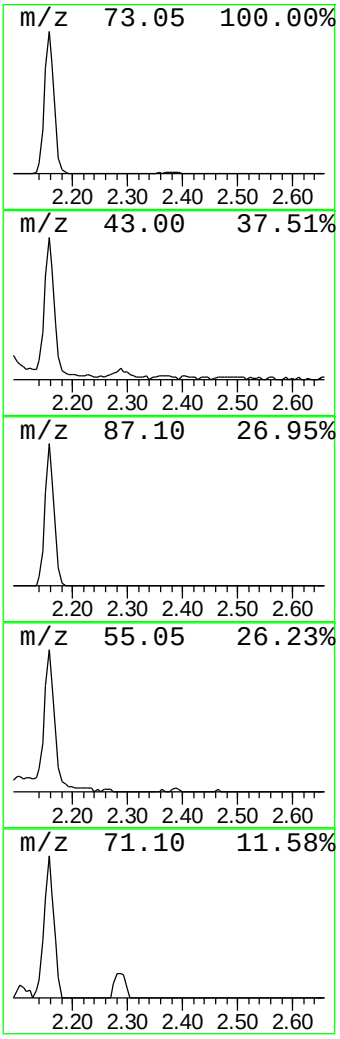
****

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.16	5.84 ng	229570	1,4-Dichlorobenzene-d4	6.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3	3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	25
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12
5	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
Data File : BF099186.D
Acq On : 7 Oct 2017 11:26
Operator : SJ/JU
Sample : I5586-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G72-0-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

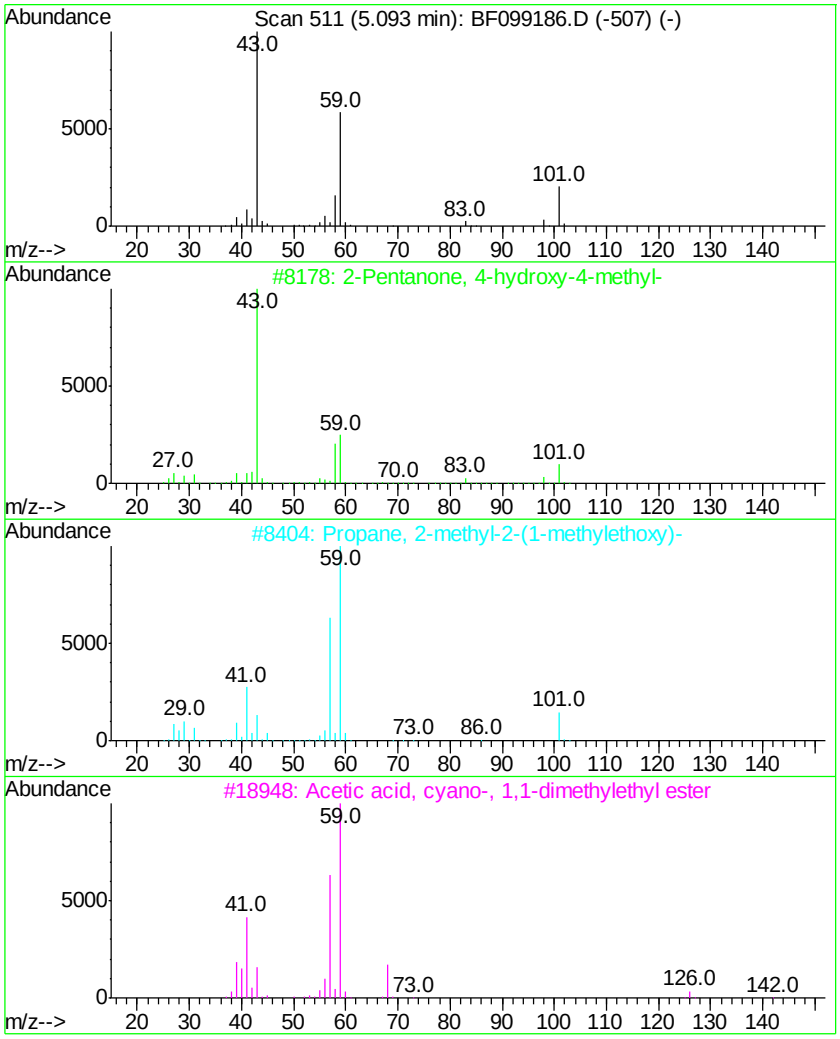
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

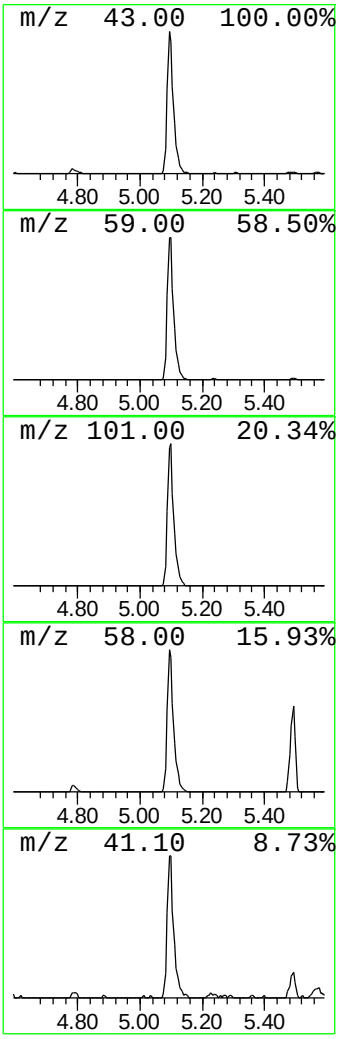
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.09	16.16 ng	635309	1,4-Dichlorobenzene-d4	6.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	28
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	17
4	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9





Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
Data File : BF099186.D
Acq On : 7 Oct 2017 11:26
Operator : SJ/JU
Sample : I5586-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G72-0-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

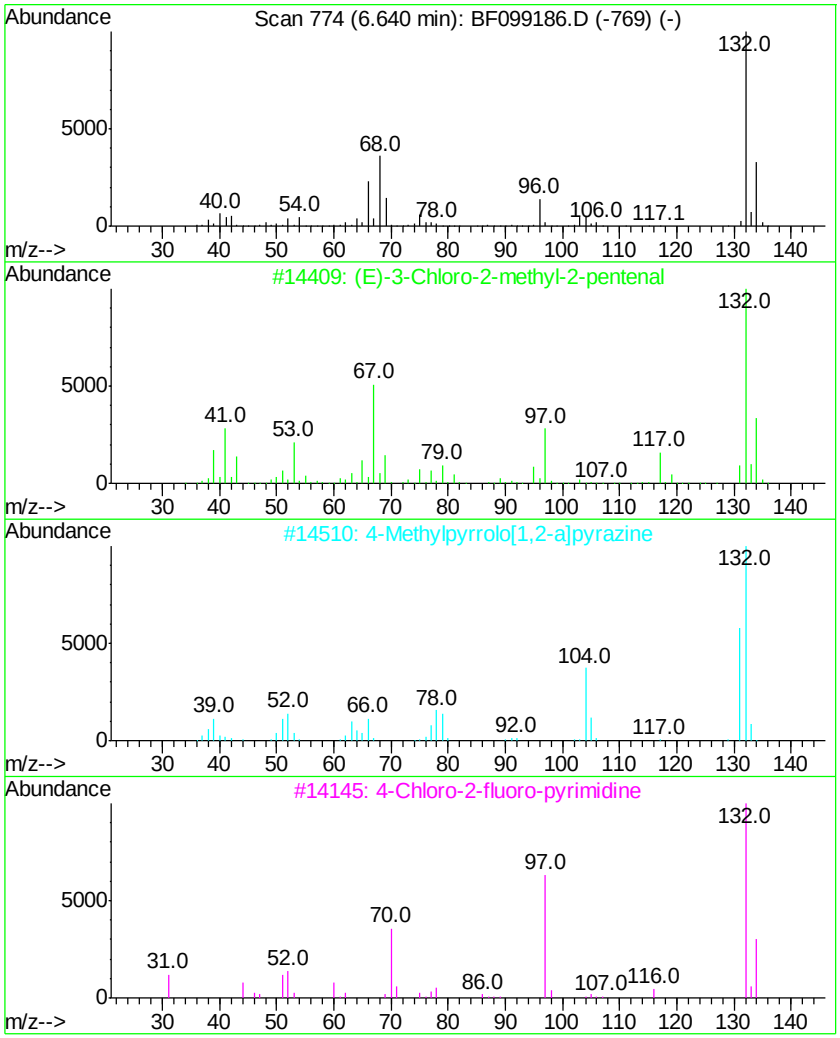
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

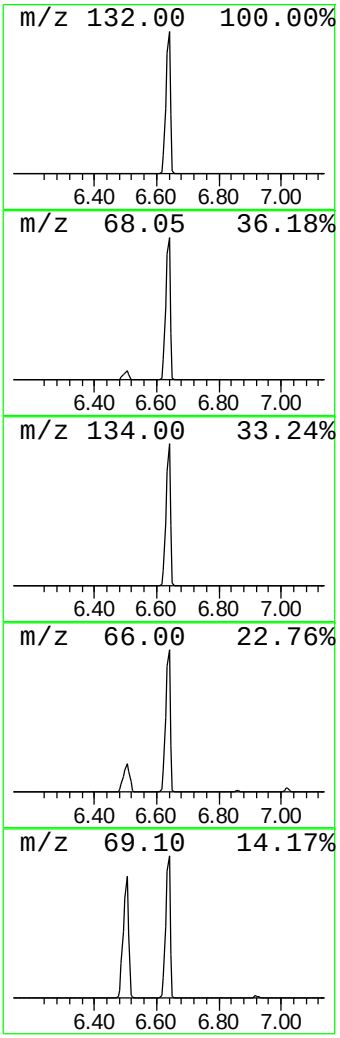
****

Peak Number 3 unknown6.64 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.64	84.29 ng	3313770	1,4-Dichlorobenzene-d4	6.86

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
2	4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9
3	4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9
4	2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9
5	Xenon	132	Xe	007440-63-3	9





Library Search Compound Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
Data File : BF099186.D
Acq On : 7 Oct 2017 11:26
Operator : SJ/JU
Sample : I5586-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G72-0-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

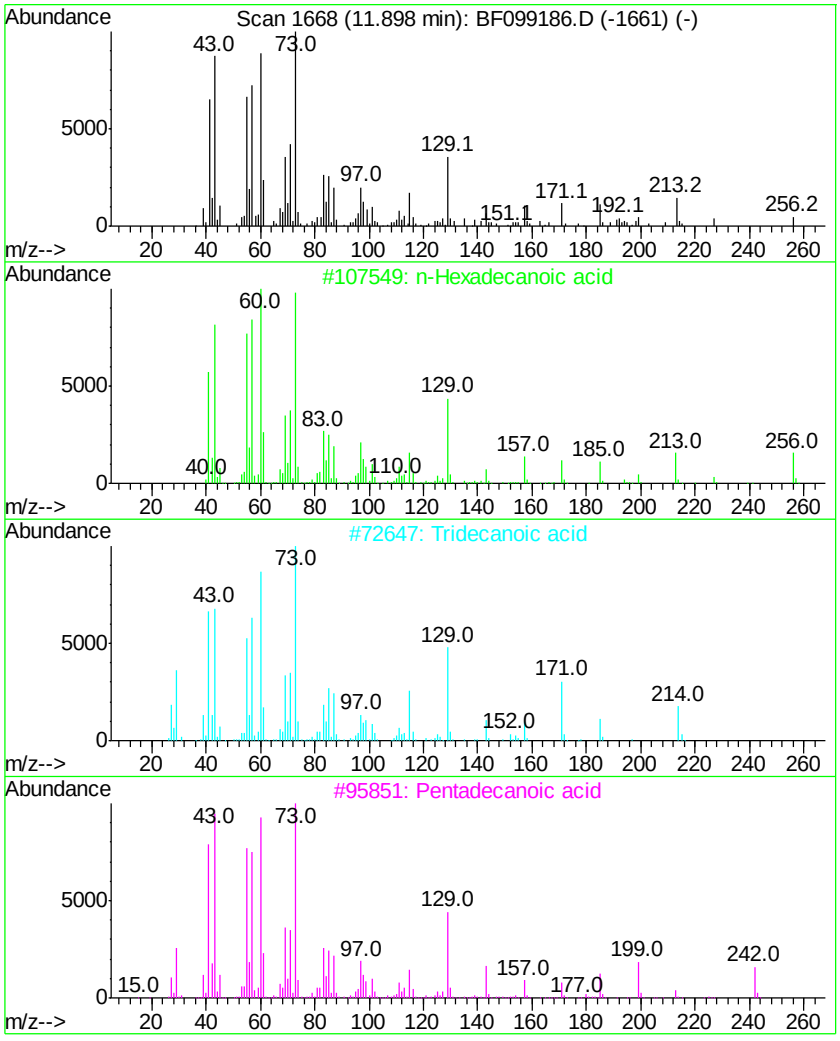
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

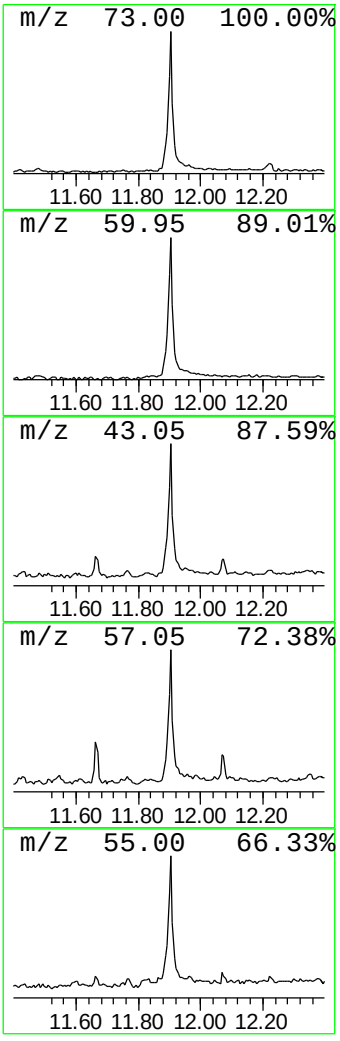
****

Peak Number 5 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.90	9.47 ng	504636	Phenanthrene-d10	11.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tridecanoic acid	214	C13H26O2	000638-53-9	95
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	93
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	91
5	Dodecanoic acid	200	C12H24O2	000143-07-7	76





Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
Data File : BF099186.D
Acq On : 7 Oct 2017 11:26
Operator : SJ/JU
Sample : I5586-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G72-0-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

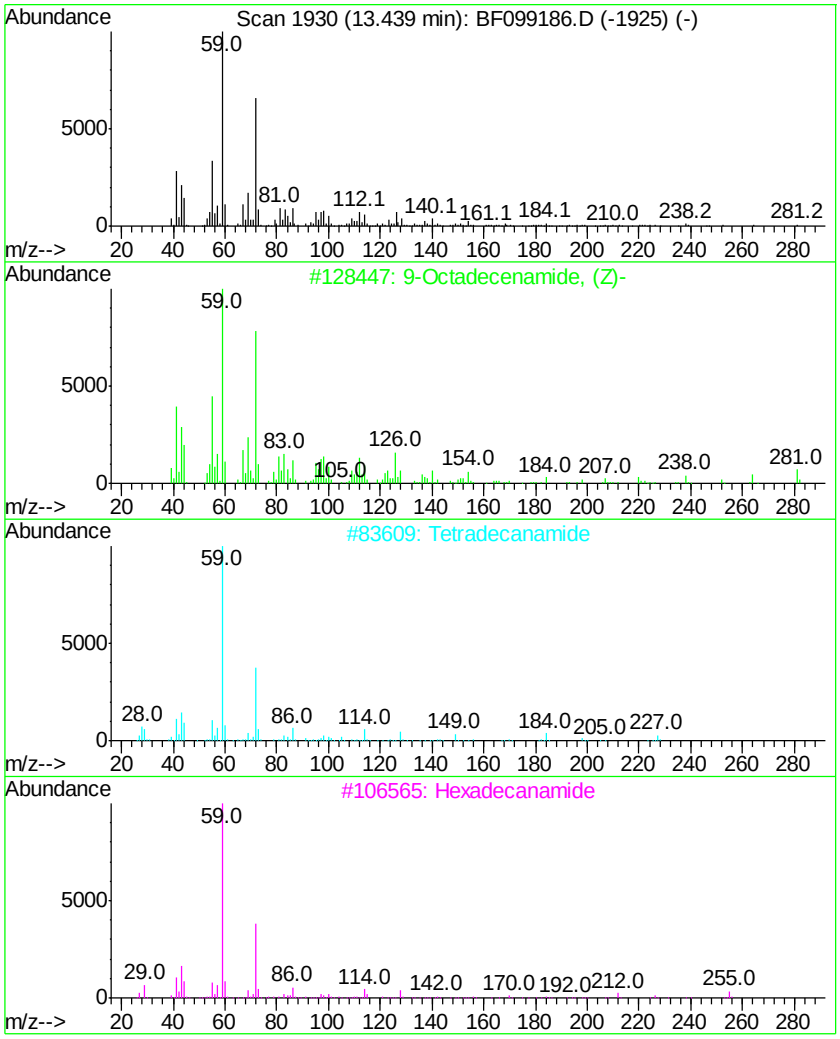
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

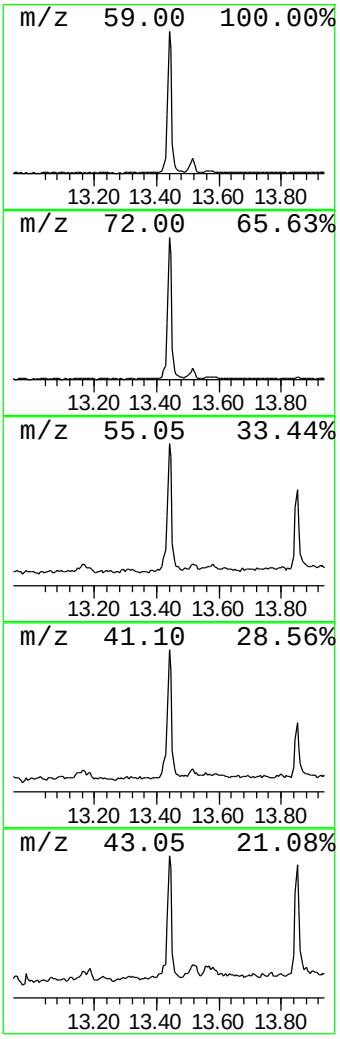
****

Peak Number 6 9-Octadecenamide, (Z)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.44	17.84 ng	691695	Chrysene-d12	14.02

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	97
2	Tetradecanamide	227	C14H29NO	000638-58-4	64
3	Hexadecanamide	255	C16H33NO	000629-54-9	59
4	Pentadecanamide, 15-bromo-	319	C15H30BrNO	1000163-86-1	59
5	Nonanamide	157	C9H19NO	001120-07-6	53





Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
Data File : BF099186.D
Acq On : 7 Oct 2017 11:26
Operator : SJ/JU
Sample : I5586-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G72-0-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

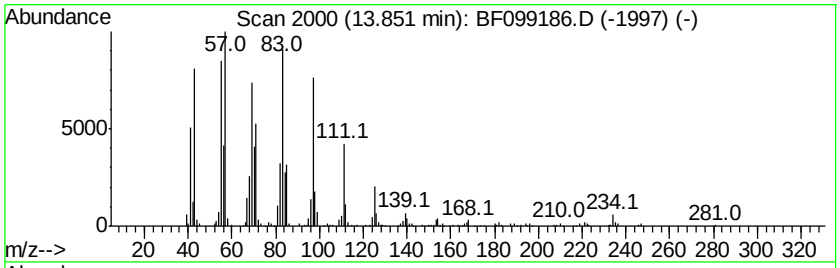
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

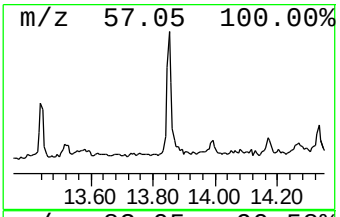
****

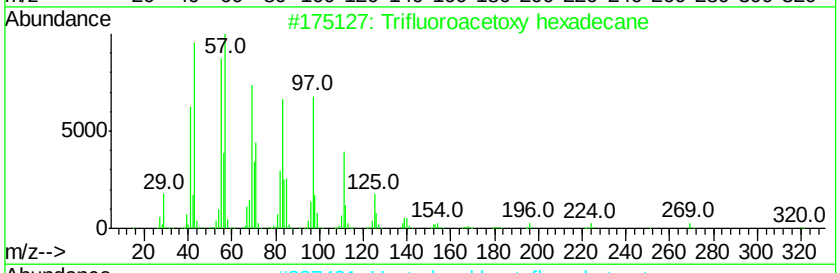
Peak Number 7 Trifluoroacetoxy hexadecane Concentration Rank 8

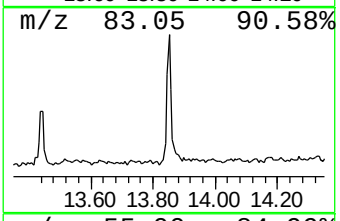
R.T.	EstConc	Area	Relative to ISTD	R.T.
13.85	8.23 ng	319176	Chrysene-d12	14.02

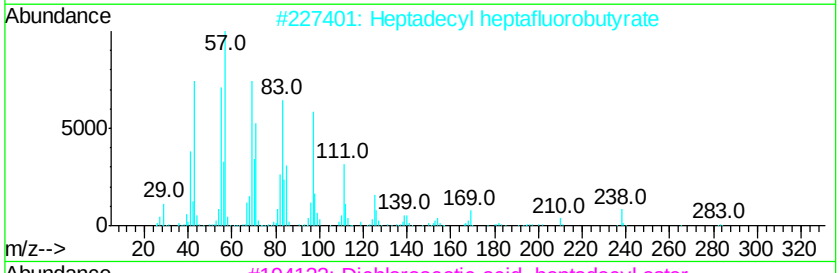
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	94
2	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
3	Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	94
4	1-Heneicosanol	312	C21H44O	015594-90-8	93
5	Cyclopentadecane	210	C15H30	000295-48-7	93

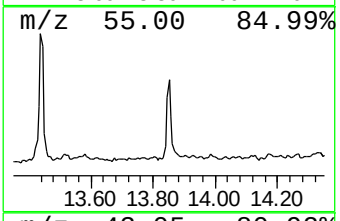


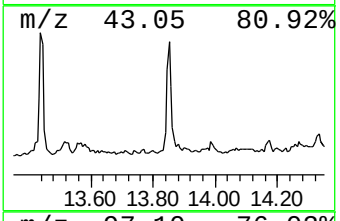


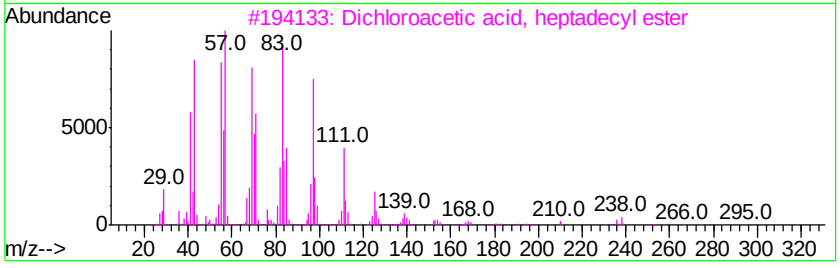


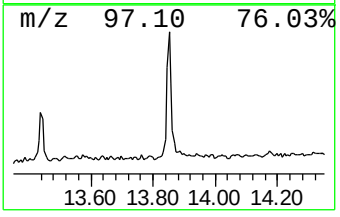












Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
Data File : BF099186.D
Acq On : 7 Oct 2017 11:26
Operator : SJ/JU
Sample : I5586-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G72-0-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

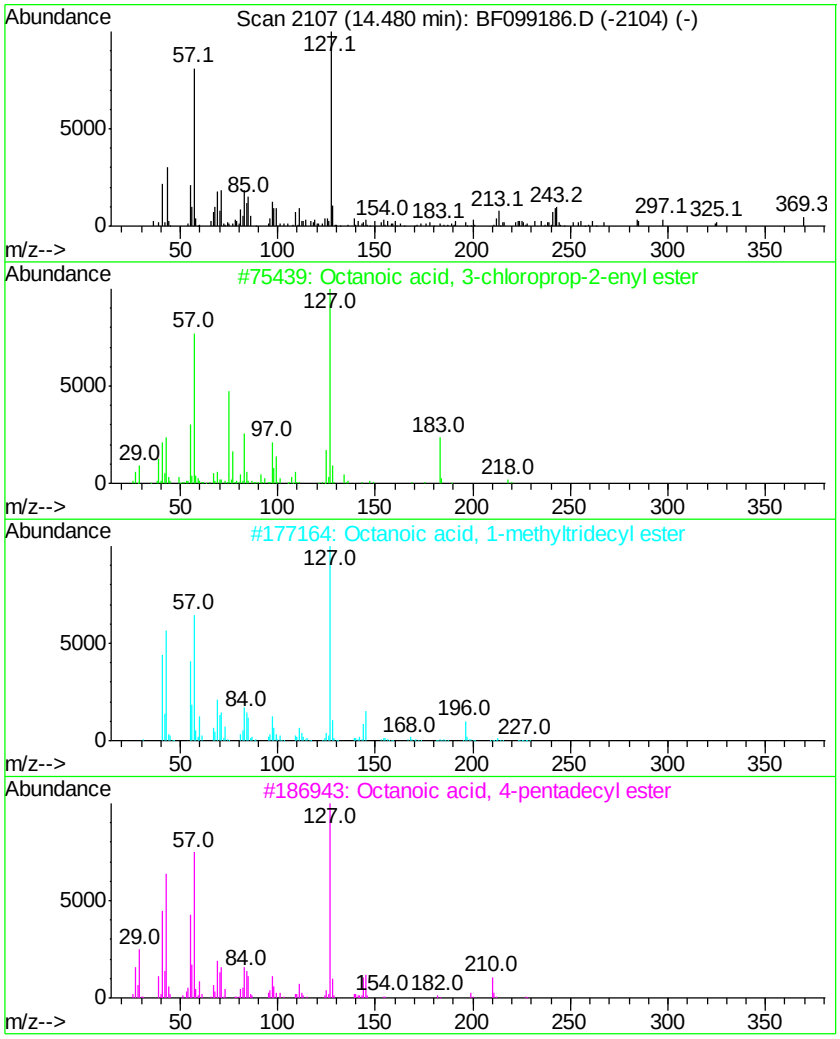
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

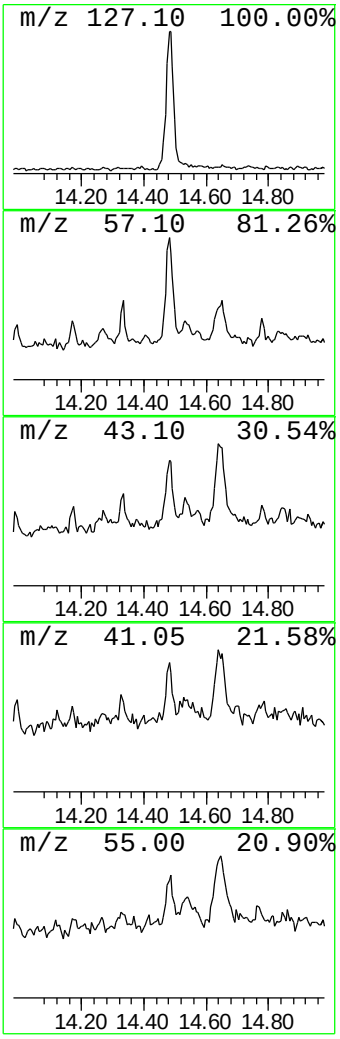
****

Peak Number 8 Octanoic acid, 3-chloroprop... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.48	3.70 ng	143425	Chrysene-d12	14.02

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octanoic acid, 3-chloroprop-2-en...	218	C11H19ClO2	1000299-35-0	53
2	Octanoic acid, 1-methyltridecyl ...	340	C22H44O2	055193-79-8	53
3	Octanoic acid, 4-pentadecyl ester	354	C23H46O2	1000280-63-2	53
4	2-Hexene, 1,1-diethoxy-	172	C10H20O2	054306-00-2	53
5	2-Hydroxy-4-hydroxylaminopyrimidine	127	C4H5N3O2	020555-88-8	52





Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
Data File : BF099186.D
Acq On : 7 Oct 2017 11:26
Operator : SJ/JU
Sample : I5586-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G72-0-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

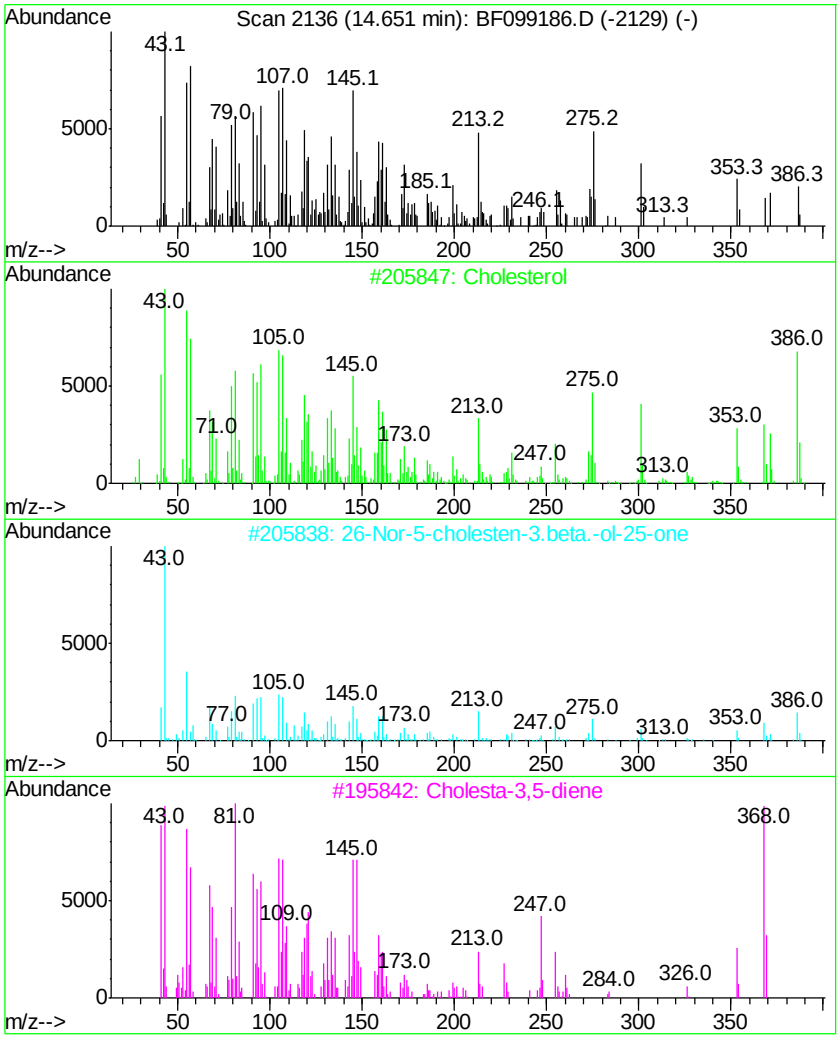
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

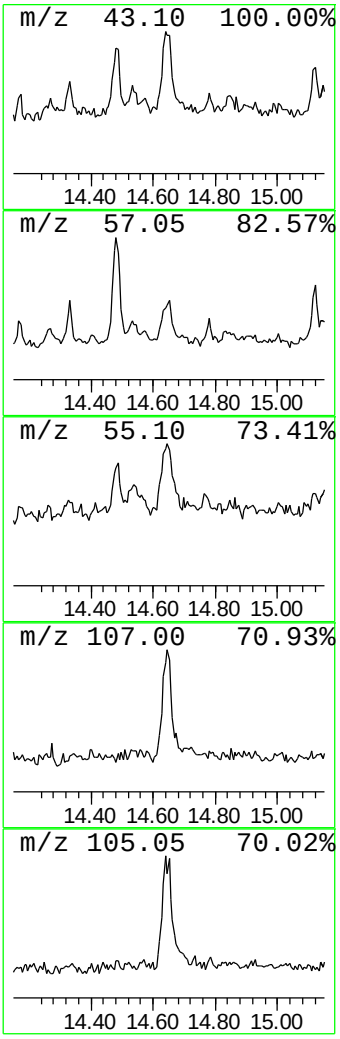
****

Peak Number 9 Cholesterol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.65	10.71 ng	415002	Chrysene-d12	14.02

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cholesterol	386	C27H46O	000057-88-5	99
2	26-Nor-5-cholesten-3.beta.-ol-25...	386	C26H42O2	007494-34-0	93
3	Cholesta-3,5-diene	368	C27H44	000747-90-0	38
4	Cholestane, 4,5-epoxy-, (4.alpha...	386	C27H46O	006079-19-2	25
5	Bicyclo[5.2.0]nonane, 2-methylen...	204	C15H24	242794-76-9	25





Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
Data File : BF099186.D
Acq On : 7 Oct 2017 11:26
Operator : SJ/JU
Sample : I5586-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G72-0-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

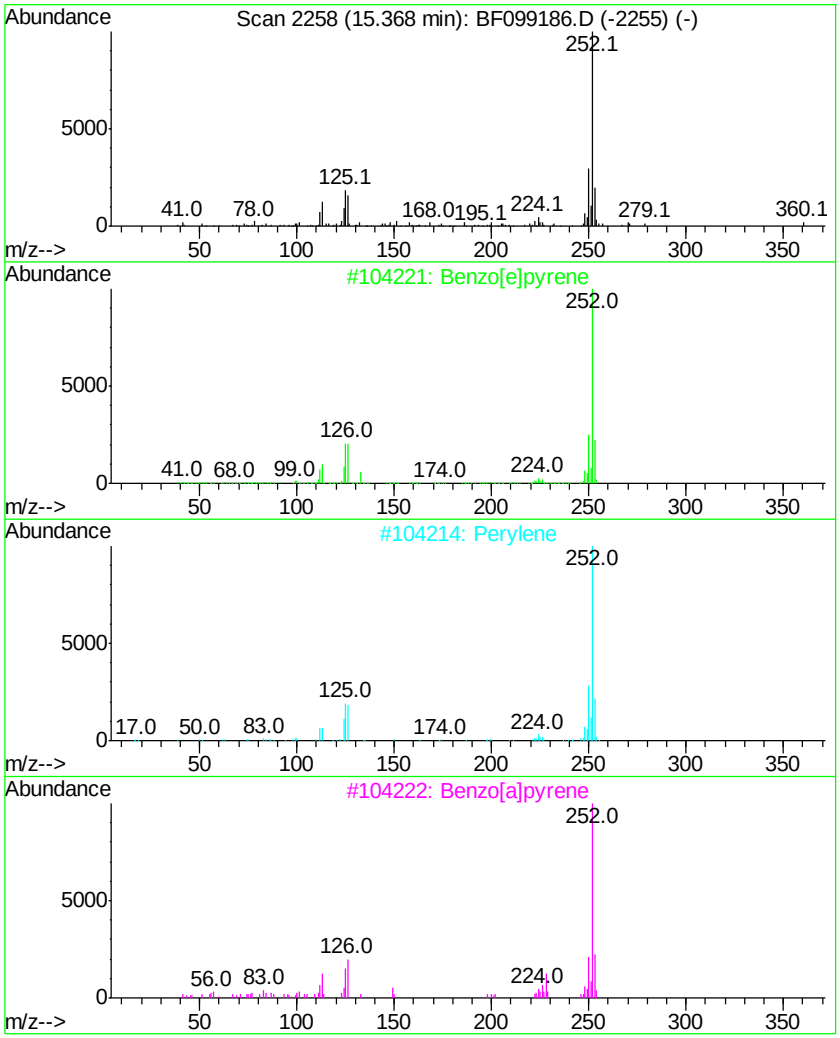
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

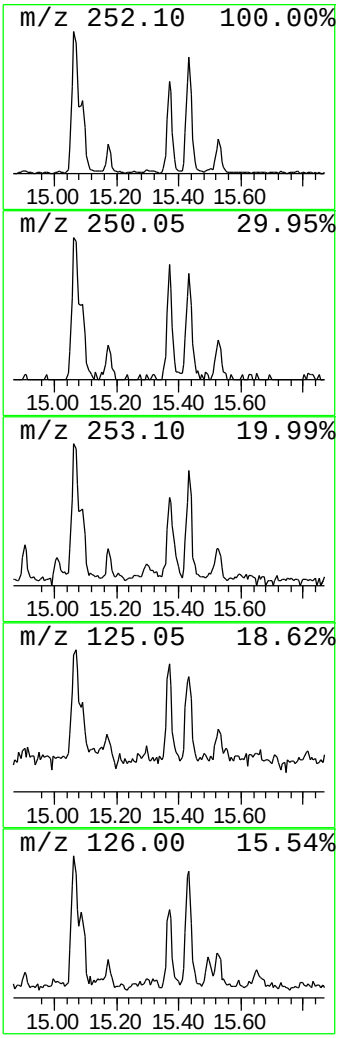
****

Peak Number 10 Benzo[e]pyrene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.37	2.23 ng	74750	Perylene-d12	15.50

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	96
2	Perylene	252	C20H12	000198-55-0	93
3	Benzo[a]pyrene	252	C20H12	000050-32-8	91
4	Benzo[k]fluoranthene	252	C20H12	000207-08-9	91
5	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	90





Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
Data File : BF099186.D
Acq On : 7 Oct 2017 11:26
Operator : SJ/JU
Sample : I5586-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G72-0-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

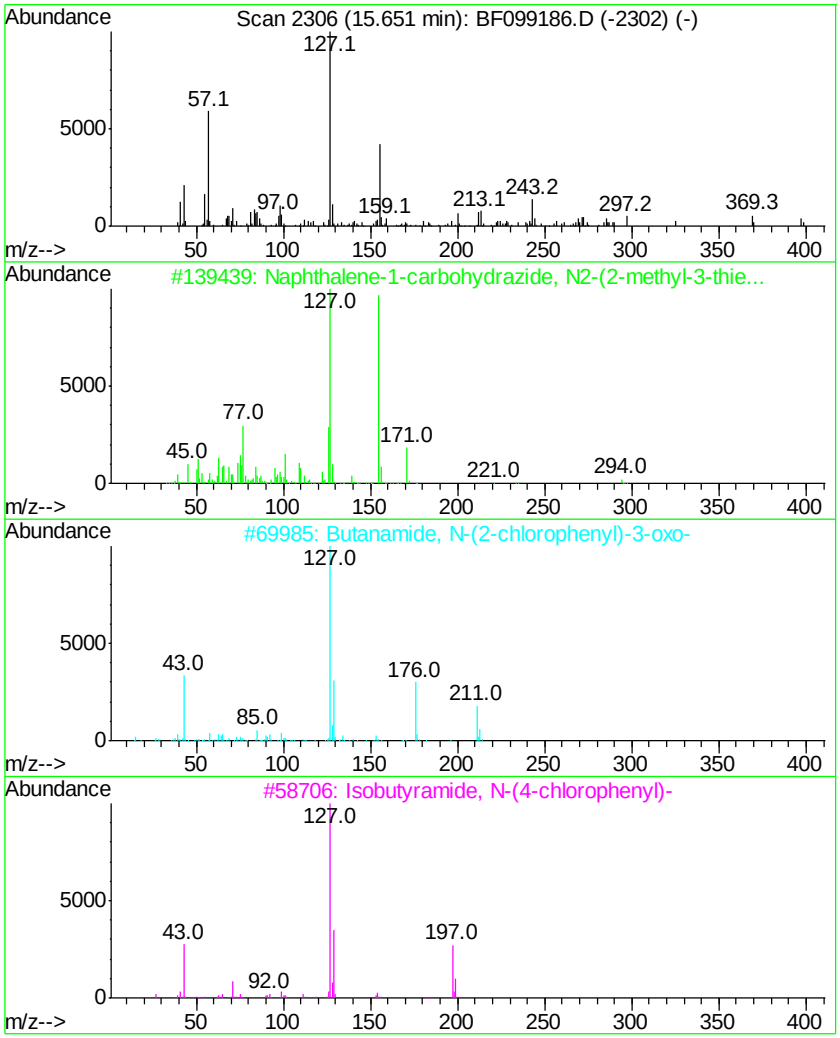
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

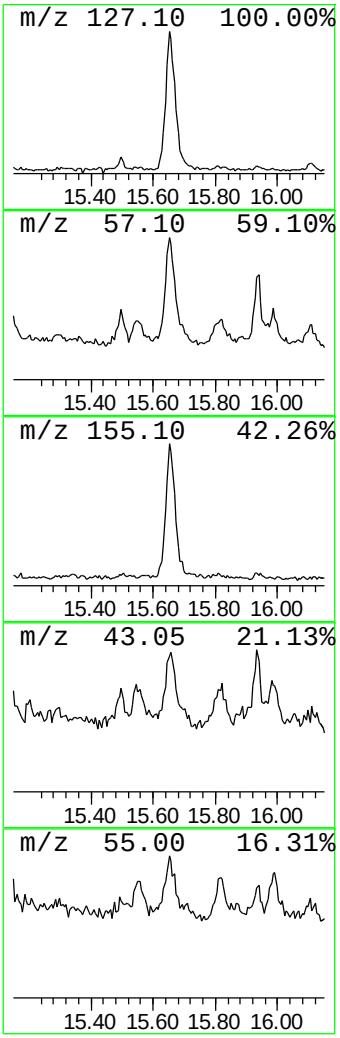
****

Peak Number 11 unknown15.65 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.65	7.06 ng	236440	Perylene-d12	15.50

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene-1-carbohydrazide, N2...	294	C17H14N2OS	1000263-63-0	45
2	Butanamide, N-(2-chlorophenyl)-3...	211	C10H10ClNO2	000093-70-9	38
3	Isobutyramide, N-(4-chlorophenyl)-	197	C10H12ClNO	1000339-96-3	35
4	Urea, N-(4-chlorophenyl)-N'-[3-(...	292	C15H17ClN2O2	1000351-08-3	35
5	Pentanamide, N-(3-chlorophenyl)-...	245	C11H13Cl2NO	1000307-36-0	35





Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
Data File : BF099186.D
Acq On : 7 Oct 2017 11:26
Operator : SJ/JU
Sample : I5586-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G72-0-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

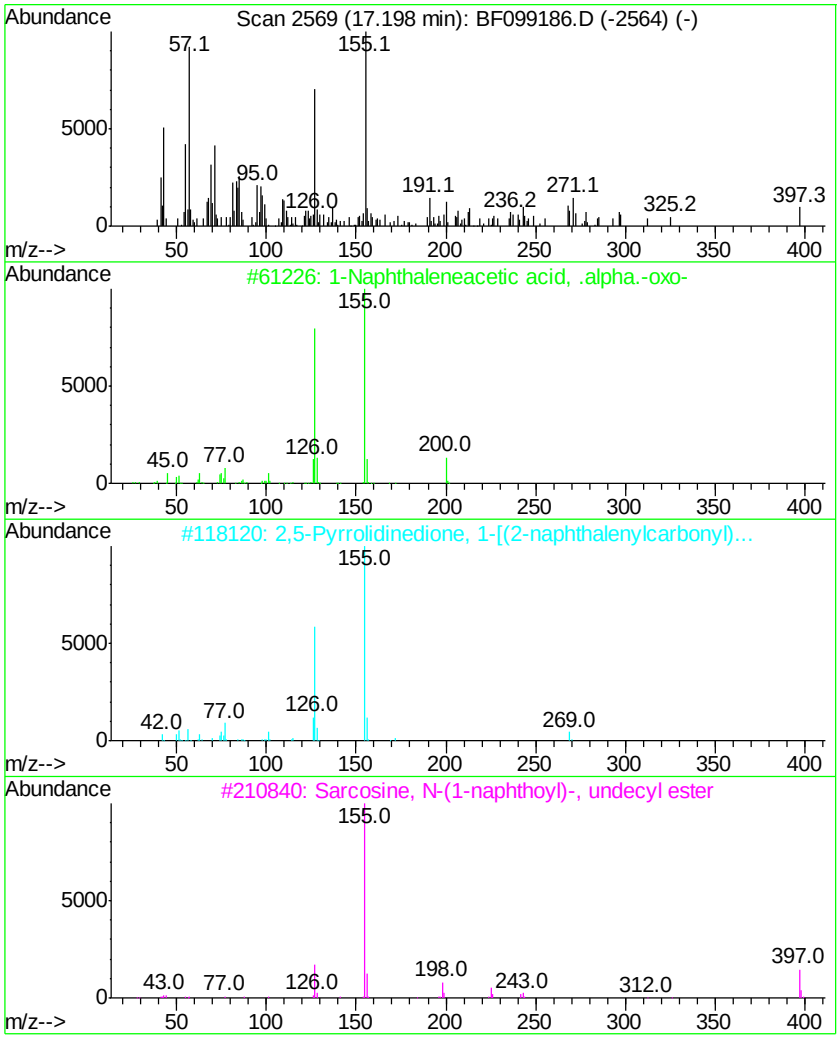
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

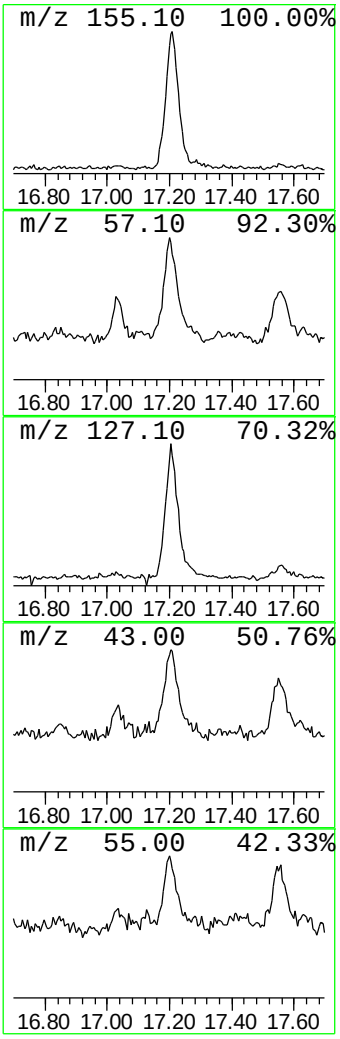
****

Peak Number 12 1-Naphthaleneacetic acid, ... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.20	8.60 ng	288163	Perylene-d12	15.50

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthaleneacetic acid, .alpha...	200	C12H8O3	026153-26-4	53
2	2,5-Pyrrolidinedione, 1-[(2-naph...	269	C15H11NO4	056374-47-1	49
3	Sarcosine, N-(1-naphthoyl)-, und...	397	C25H35NO3	1000321-40-9	43
4	5-Oxazolidinone, 2-(1,1-dimethyl...	297	C18H19NO3	119215-56-4	43
5	Fumaric acid, butyl 4-chloro3-me...	296	C15H17ClO4	1000344-86-6	38





Data Path : Z:\HPCHEM1\BNA F\DATA\BF100717\
Data File : BF099186.D
Acq On : 7 Oct 2017 11:26
Operator : SJ/JU
Sample : I5586-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G72-0-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

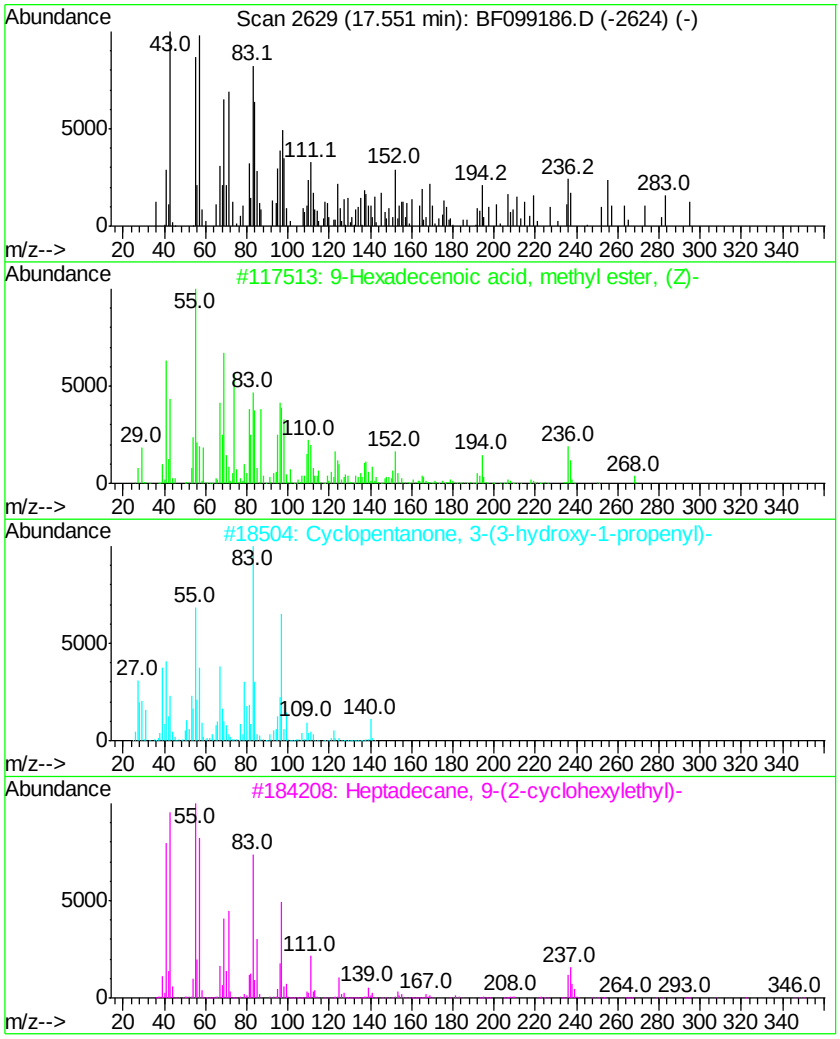
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

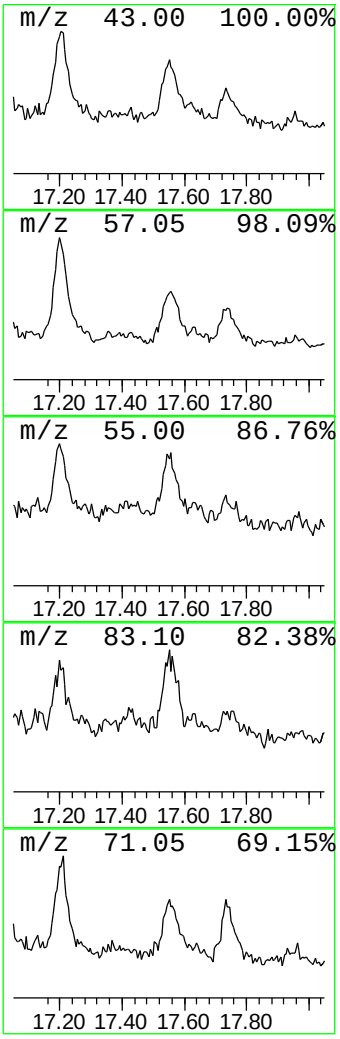
****

Peak Number 13 unknown17.55 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.55	6.49 ng	217569	Perylene-d12	15.50

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Hexadecenoic acid, methyl este...	268	C17H32O2	001120-25-8	49
2	Cyclopentanone, 3-(3-hydroxy-1-p...	140	C8H12O2	074473-08-8	43
3	Heptadecane, 9-(2-cyclohexylethyl)-	350	C25H50	025446-35-9	38
4	Cyclohexane, (3-cyclopentylpropyl)-	194	C14H26	002883-07-0	38
5	Trichloroacetic acid, dodecyl ester	330	C14H25Cl3O2	074339-50-7	38





Tentatively Identified Compound (LSC) summary

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100717\
Data File : BF099186.D
Acq On : 7 Oct 2017 11:26
Operator : SJ/JU
Sample : I5586-01
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : G72-0-1

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-methoxy...	2.16	5.8	ng	229570	1	6.86	786297	20.0
2-Pentanone, 4-hy...	5.09	16.2	ng	635309	1	6.86	786297	20.0
unknown6.64	6.64	84.3	ng	3313770	1	6.86	786297	20.0
n-Hexadecanoic acid	11.90	9.5	ng	504636	4	11.39	1065660	20.0
9-Octadecenamide,...	13.44	17.8	ng	691695	5	14.02	775267	20.0
Trifluoroacetoxy ...	13.85	8.2	ng	319176	5	14.02	775267	20.0
Octanoic acid, 3-...	14.48	3.7	ng	143425	5	14.02	775267	20.0
Cholesterol	14.65	10.7	ng	415002	5	14.02	775267	20.0
Benzo[e]pyrene	15.37	2.2	ng	74750	6	15.50	670015	20.0
unknown15.65	15.65	7.1	ng	236440	6	15.50	670015	20.0
1-Naphthaleneacet...	17.20	8.6	ng	288163	6	15.50	670015	20.0
unknown17.55	17.55	6.5	ng	217569	6	15.50	670015	20.0