LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112619\
Data File : BF118015.D
Acq On : 26 Nov 2019 18:56
Operator : JU
Sample : K5948-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : B-1-0-6

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.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	2.193	10	18	23	rVB	527626	680750	22.69%	2.450%
2	5.110	509	514	522	rVB	434796	522291	17.41%	1.880%
3	5.199	525	529	534	rBV	79372	80252	2.68%	0.289%
4	5.516	578	583	586	rBV	2754107	2592538	86.42%	9.332%
5	6.528	749	755	765	rBV	2300173	2411753	80.39%	8.681%
6	6.663	774	778	782	rBV	2594719	2494143	83.14%	8.978%
7	6.893	814	817	825	rVB	793687	650916	21.70%	2.343%
8	7.051	840	844	847	rBV	2279415	1952058	65.07%	7.027%
9	7.457	908	913	916	rBV	1653323	1466614	48.89%	5.279%
10	8.181	1030	1036	1039	rBV	1007298	832508	27.75%	2.997%
11	8.204	1039	1040	1043	rVB	75917	36407	1.21%	0.131%
12	9.263	1215	1220	1223	rBV	3242794	3000040	100.00%	10.799%
13	9.339	1230	1233	1235	rBV	61828	53526	1.78%	0.193%
14	9.375	1235	1239	1241	rVB2	28968	32796	1.09%	0.118%
15	9.651	1283	1286	1290	rVB	552788	425530	14.18%	1.532%
16	9.804	1310	1312	1315	rBV2	36084	40584	1.35%	0.146%
17	9.869	1319	1323	1326	rVV2	58288	63483	2.12%	0.229%
18	9.945	1328	1336	1339	rVV	1188839	973199	32.44%	3.503%
19	9.981	1339	1342	1350	rVV5	40065	61931	2.06%	0.223%
20	10.045	1350	1353	1358	rVV4	22136	32957	1.10%	0.119%
21	10.145	1364	1370	1372	rBV6	36029	46307	1.54%	0.167%
22	10.263	1387	1390	1392	rBV2	29922	34466	1.15%	0.124%
23	10.375	1407	1409	1411	rVB	54977	41490	1.38%	0.149%
24	10.592	1442	1446	1451	rBV3	33035	43220	1.44%	0.156%
25	10.739	1467	1471	1475	rVB	1694976	1559360	51.98%	5.613%
26	10.863	1486	1492	1498	rVB3	89427	149065	4.97%	0.537%
27	11.075	1525	1528	1530	rBV2	36485	35191	1.17%	0.127%
28	11.298	1563	1566	1568	rBV	59112	54541	1.82%	0.196%
29	11.333	1569	1572	1576	rVB2	53331	65843	2.19%	0.237%
30	11.439	1586	1590	1592	rBV	837346	698899	23.30%	2.516%
31	11.463	1592	1594	1597	rVV	136664	109332	3.64%	0.394%
32	11.510	1599	1602	1605	rVB2	44130	50830	1.69%	0.183%
33	11.669	1627	1629	1635	rBV2	37965	51912	1.73%	0.187%
34	11.728	1636	1639	1642	rVB2	43435	42216	1.41%	0.152%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112619\
Data File : BF118015.D
Acq On : 26 Nov 2019 18:56
Operator : JU
Sample : K5948-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : B-1-0-6

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	11.963	1674	1679	1683	rBV2	358946	413447	13.78%	1.488%
36	12.004	1684	1686	1692	rVB4	68332	86902	2.90%	0.313%
37	12.051	1692	1694	1697	rBV2	63956	59195	1.97%	0.213%
38	12.492	1767	1769	1772	rBV	46514	59632	1.99%	0.215%
39	12.522	1772	1774	1778	rVB3	91049	104699	3.49%	0.377%
40	12.657	1794	1797	1801	rBV	209316	191140	6.37%	0.688%
41	12.698	1801	1804	1809	rVB4	37484	39443	1.31%	0.142%
42	12.745	1809	1812	1816	rBV2	138345	142130	4.74%	0.512%
43	12.886	1831	1836	1842	rBV	237041	355785	11.86%	1.281%
44	13.027	1856	1860	1864	rBV	2514863	2228364	74.28%	8.021%
45	13.216	1889	1892	1894	rBV	60754	59177	1.97%	0.213%
46	13.322	1908	1910	1914	rVB	45832	41353	1.38%	0.149%
47	13.445	1929	1931	1938	rVB3	39146	51998	1.73%	0.187%
48	13.722	1976	1978	1981	rBV3	41536	43636	1.45%	0.157%
49	13.786	1987	1989	1992	rBV3	38288	45547	1.52%	0.164%
50	13.933	2011	2014	2019	rVB4	134044	199013	6.63%	0.716%
51	14.063	2034	2036	2037	rBV	62812	55127	1.84%	0.198%
52	14.086	2037	2040	2043	rVV2	806376	833123	27.77%	2.999%
53	14.116	2043	2045	2047	rVB	134122	115289	3.84%	0.415%
54	14.551	2116	2119	2123	rBV2	129798	184102	6.14%	0.663%
55	14.845	2167	2169	2177	rVB3	143419	259310	8.64%	0.933%
56	15.592	2292	2296	2307	rVB2	441558	825143	27.50%	2.970%

Sum of corrected areas: 27780503

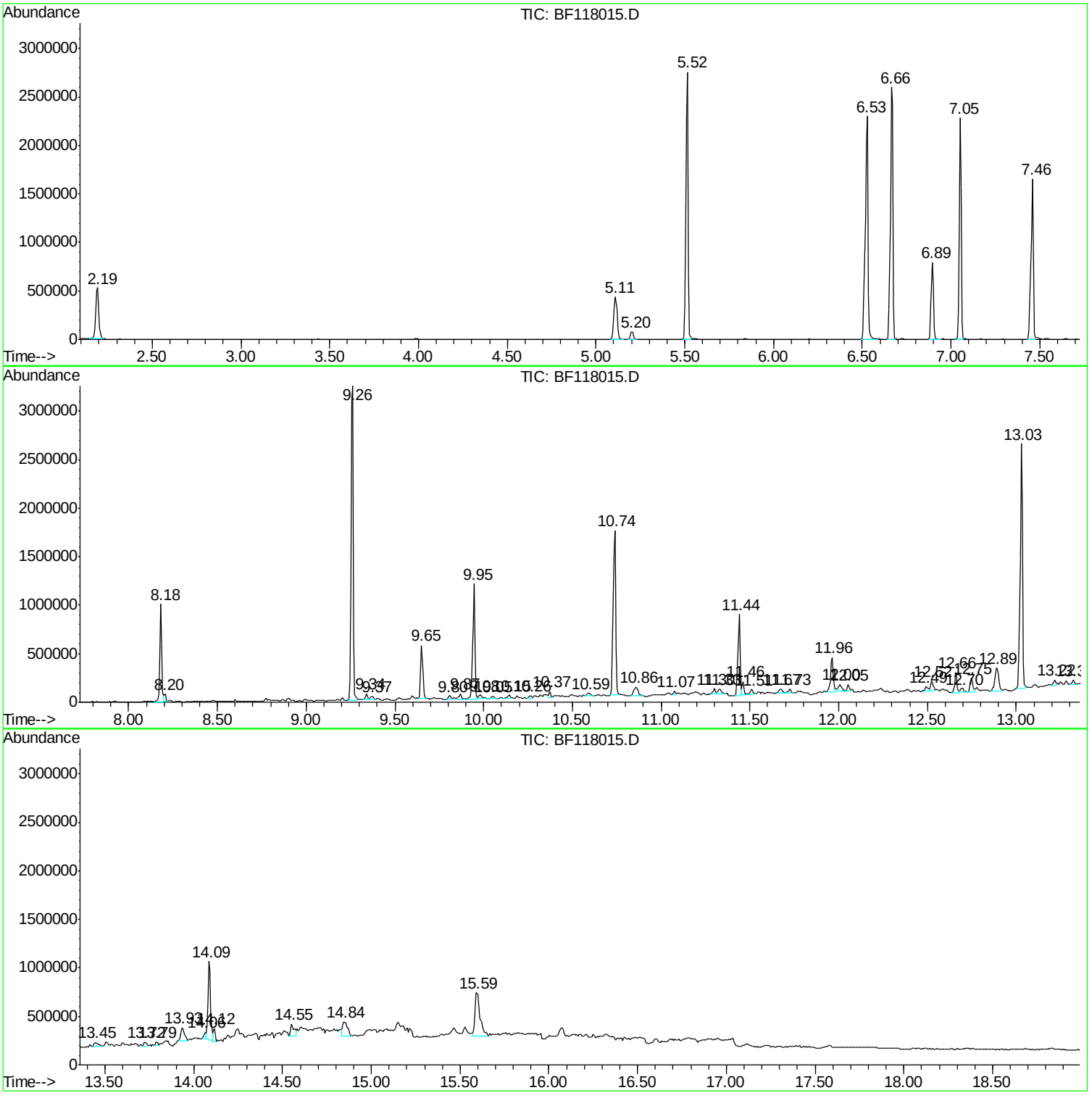
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112619\
Data File : BF118015.D
Acq On : 26 Nov 2019 18:56
Operator : JU
Sample : K5948-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : B-1-0-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112619\
Data File : BF118015.D
Acq On : 26 Nov 2019 18:56
Operator : JU
Sample : K5948-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : B-1-0-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

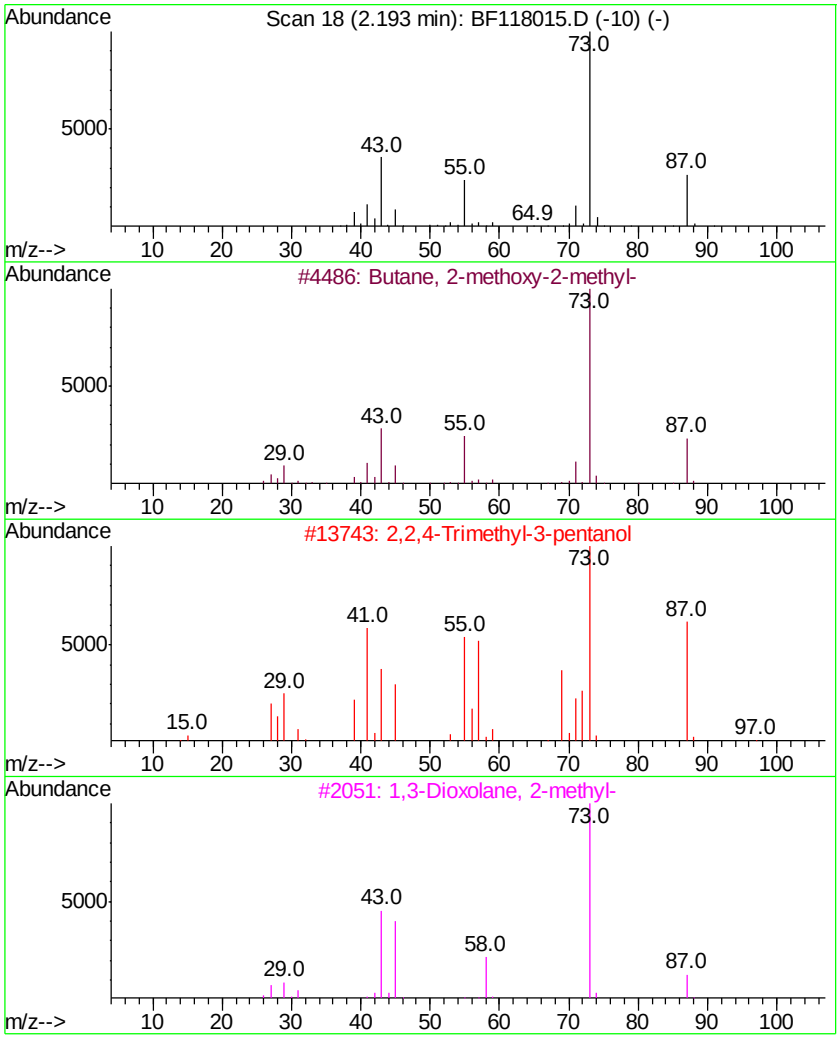
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

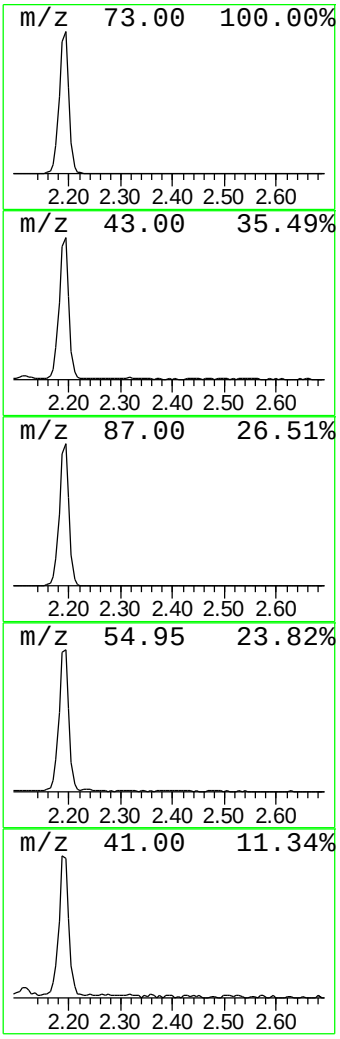
****

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.19	20.92 ng	680750	1,4-Dichlorobenzene-d4	6.89

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	56
3	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
5	Silane, tetramethyl-	88	C4H12Si	000075-76-3	17





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112619\
Data File : BF118015.D
Acq On : 26 Nov 2019 18:56
Operator : JU
Sample : K5948-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : B-1-0-6

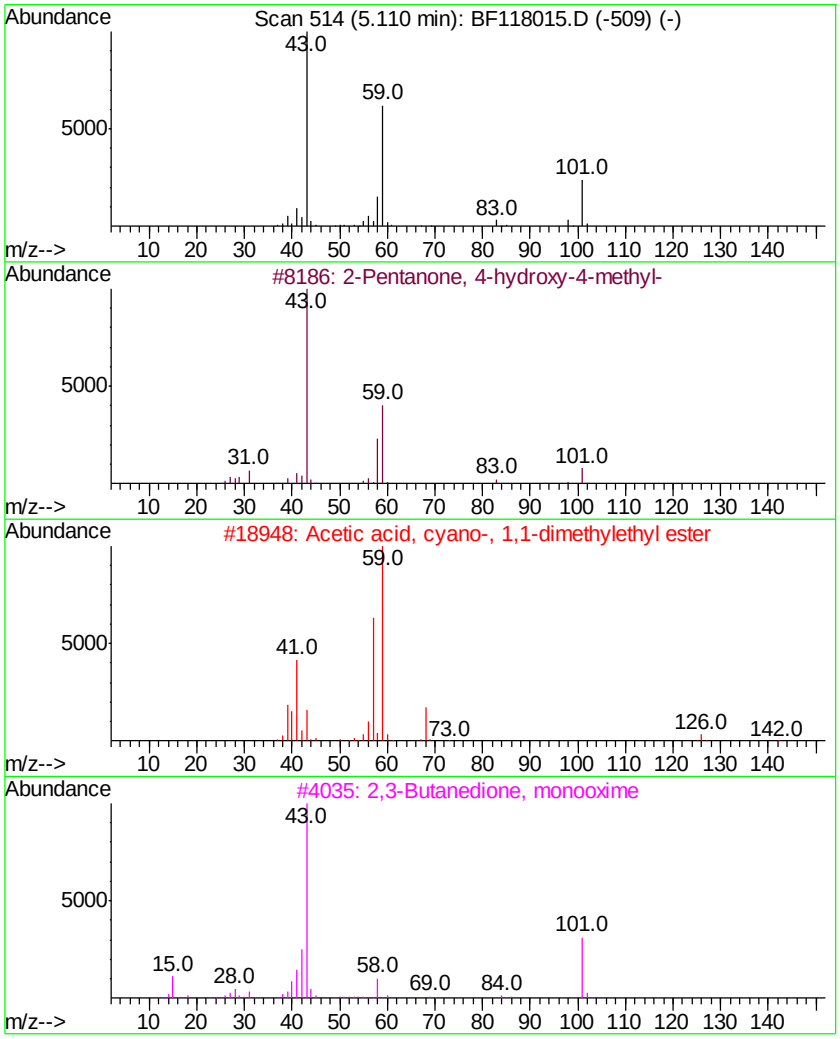
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

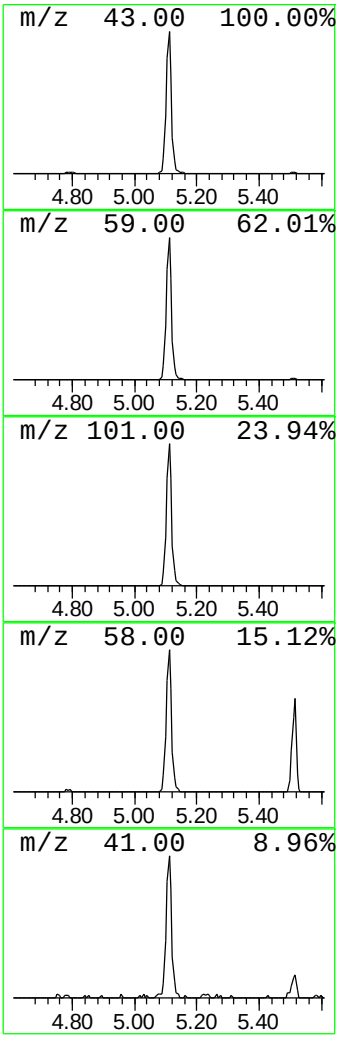
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.11	16.05 ng	522291	1,4-Dichlorobenzene-d4	6.89

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	25
3	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	17
4	(+-)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112619\
Data File : BF118015.D
Acq On : 26 Nov 2019 18:56
Operator : JU
Sample : K5948-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : B-1-0-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

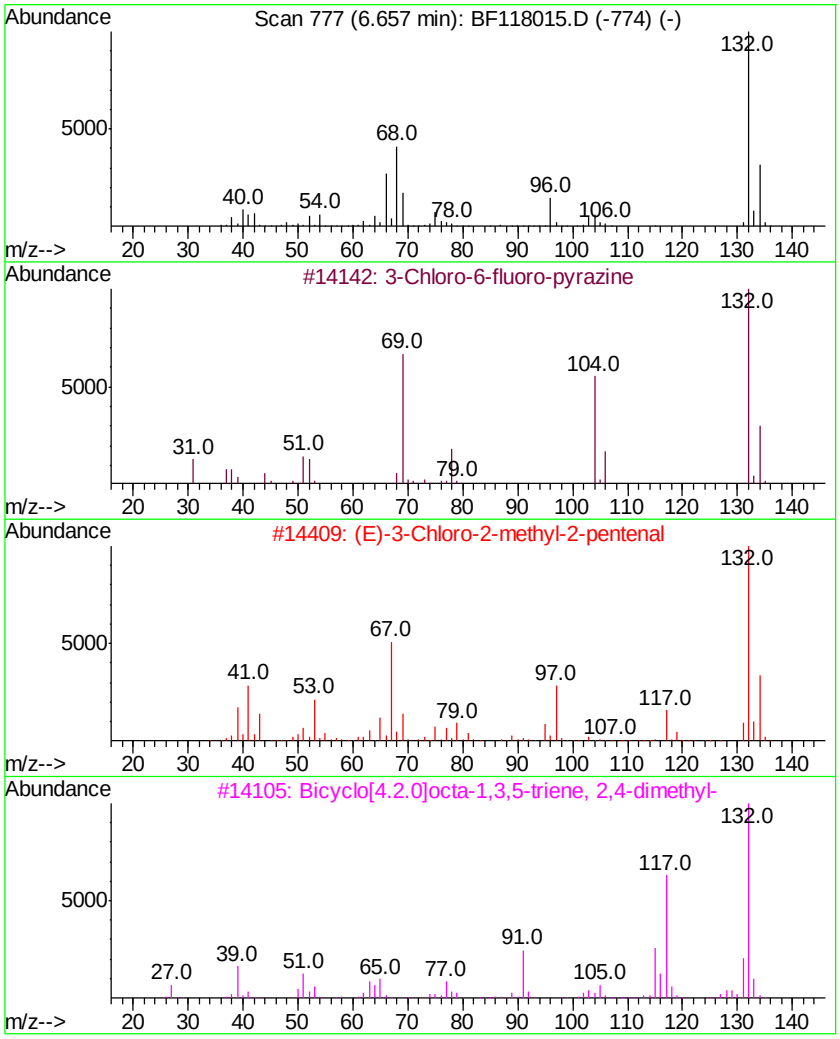
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

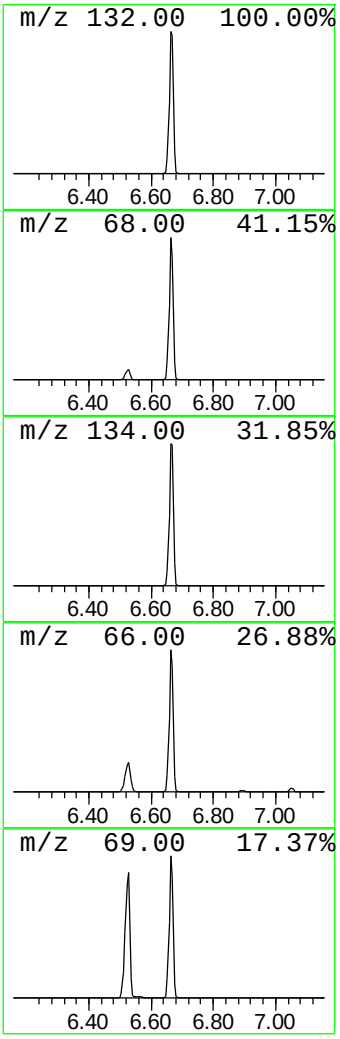
****

Peak Number 4 unknown6.66 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.66	76.63 ng	2494140	1,4-Dichlorobenzene-d4	6.89

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
4	Benzene, 1-ethenyl-3,5-dimethyl-	132	C10H12	005379-20-4	9
5	1,3,4-Thiadiazole-2(3H)-thione, ...	132	C3H4N2S2	029490-19-5	9





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112619\
Data File : BF118015.D
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Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : B-1-0-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

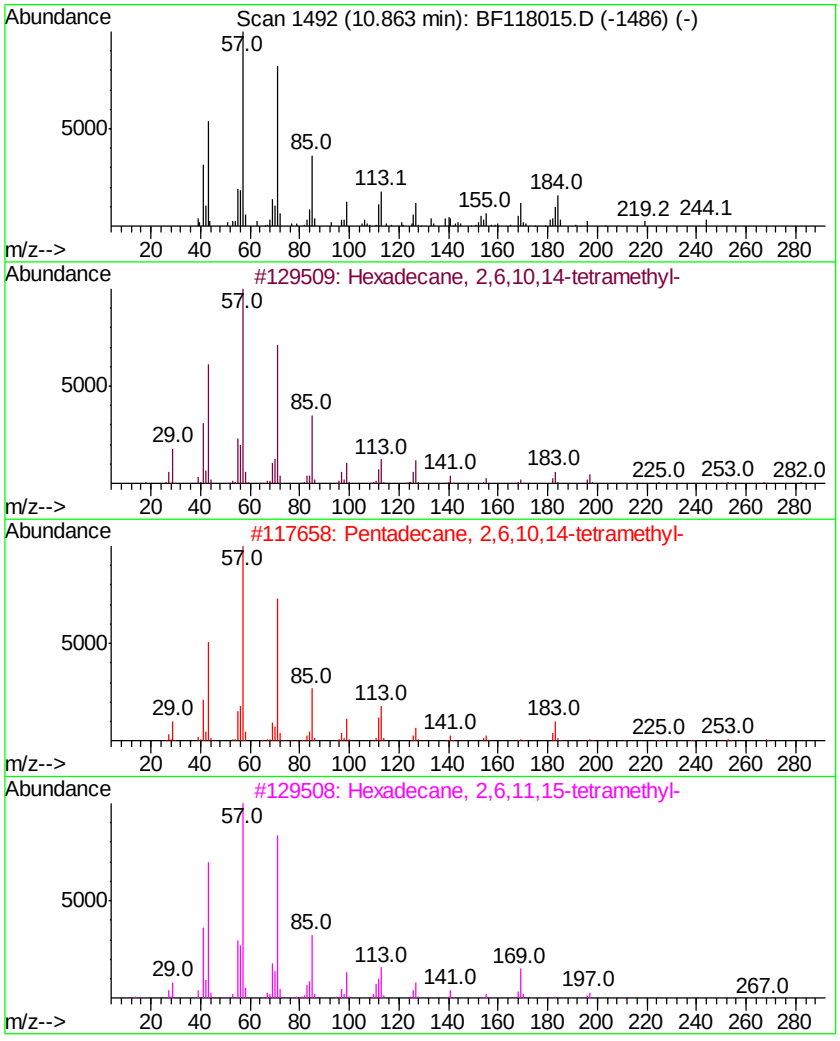
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

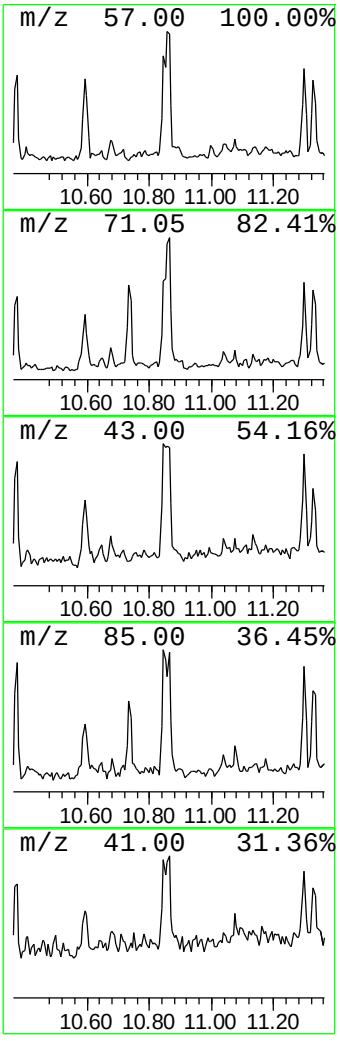
****

Peak Number 6 Hexadecane, 2,6,10,14-tetra... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.86	4.27 ng	149065	Phenanthrene-d10	11.44

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	87
2	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	86
3	Hexadecane, 2,6,11,15-tetramethyl-	282	C20H42	000504-44-9	72
4	Dodecane	170	C12H26	000112-40-3	68
5	Heptadecane, 7-methyl-	254	C18H38	020959-33-5	64





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Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : B-1-0-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

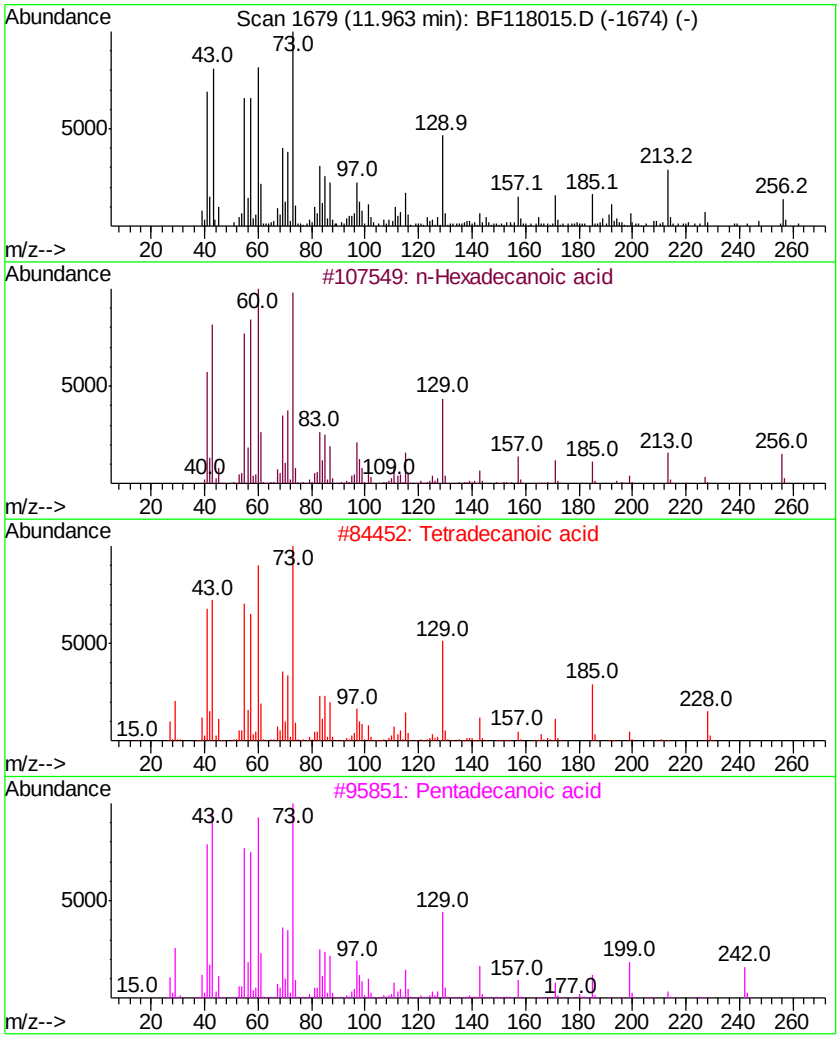
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

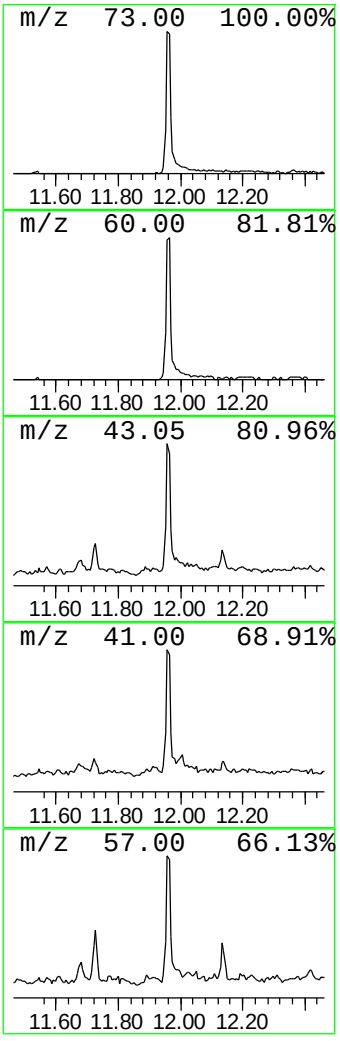
****

Peak Number 7 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.96	11.83 ng	413447	Phenanthrene-d10	11.44

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	91
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	87
4	Tridecanoic acid	214	C13H26O2	000638-53-9	86
5	n-Decanoic acid	172	C10H20O2	000334-48-5	83





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Sample : K5948-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : B-1-0-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

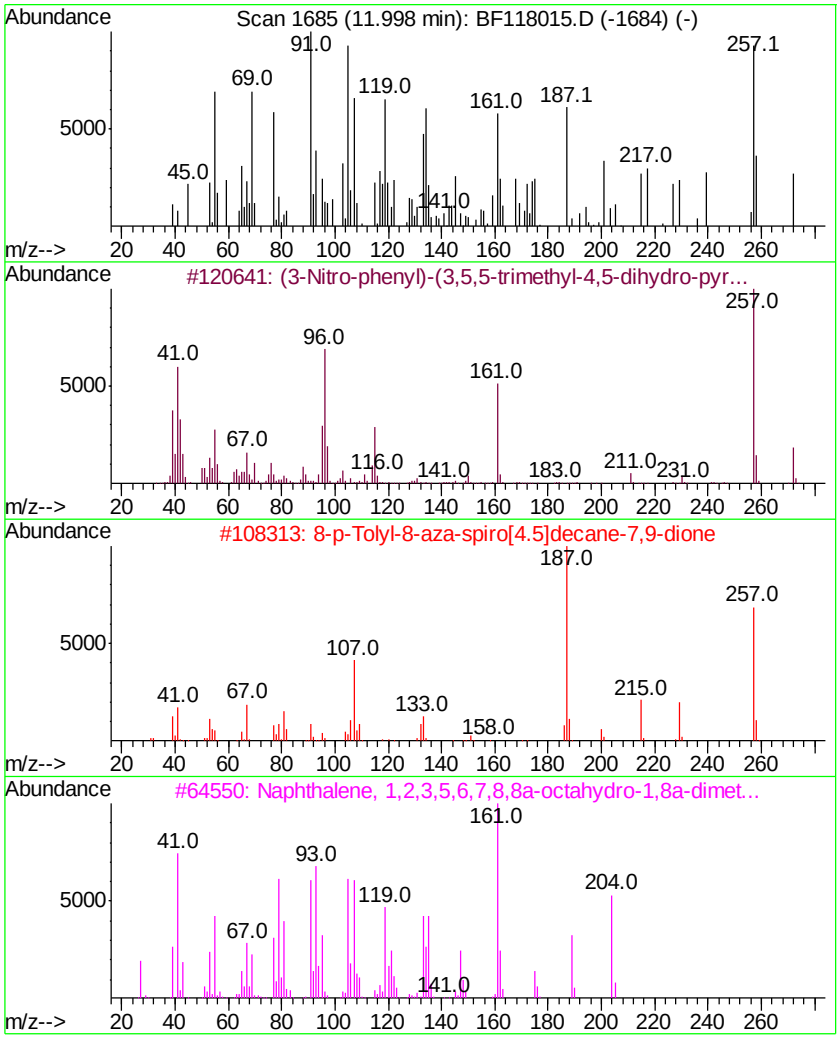
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

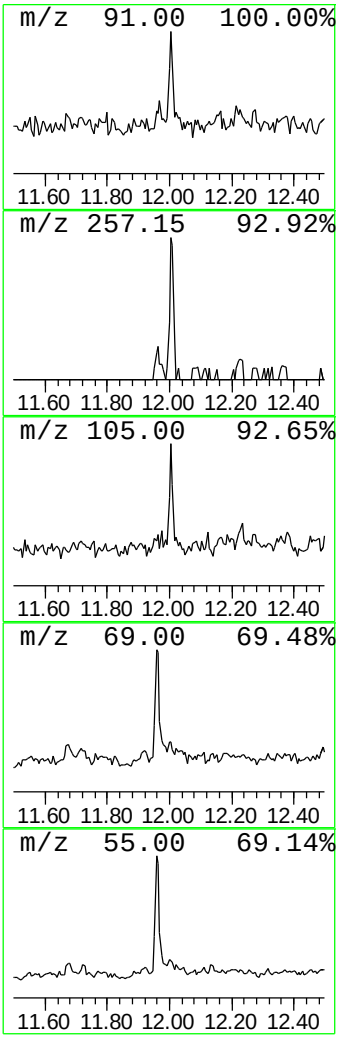
****

Peak Number 8 unknown12.00 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.00	2.49 ng	86902	Phenanthrene-d10	11.44

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(3-Nitro-phenyl)-(3,5,5-trimethy...	272	C14H16N4O2	1000274-25-3	38
2	8-p-Tolyl-8-aza-spiro[4.5]decane...	257	C16H19NO2	1000318-33-1	30
3	Naphthalene, 1,2,3,5,6,7,8,8a-oc...	204	C15H24	010219-75-7	25
4	Tricyclo[4.1.0.0(2,4)]heptane, 3...	204	C15H24	056348-21-1	25
5	1,6-Cyclodecadiene, 1-methyl-5-m...	204	C15H24	023986-74-5	25





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112619\
Data File : BF118015.D
Acq On : 26 Nov 2019 18:56
Operator : JU
Sample : K5948-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : B-1-0-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

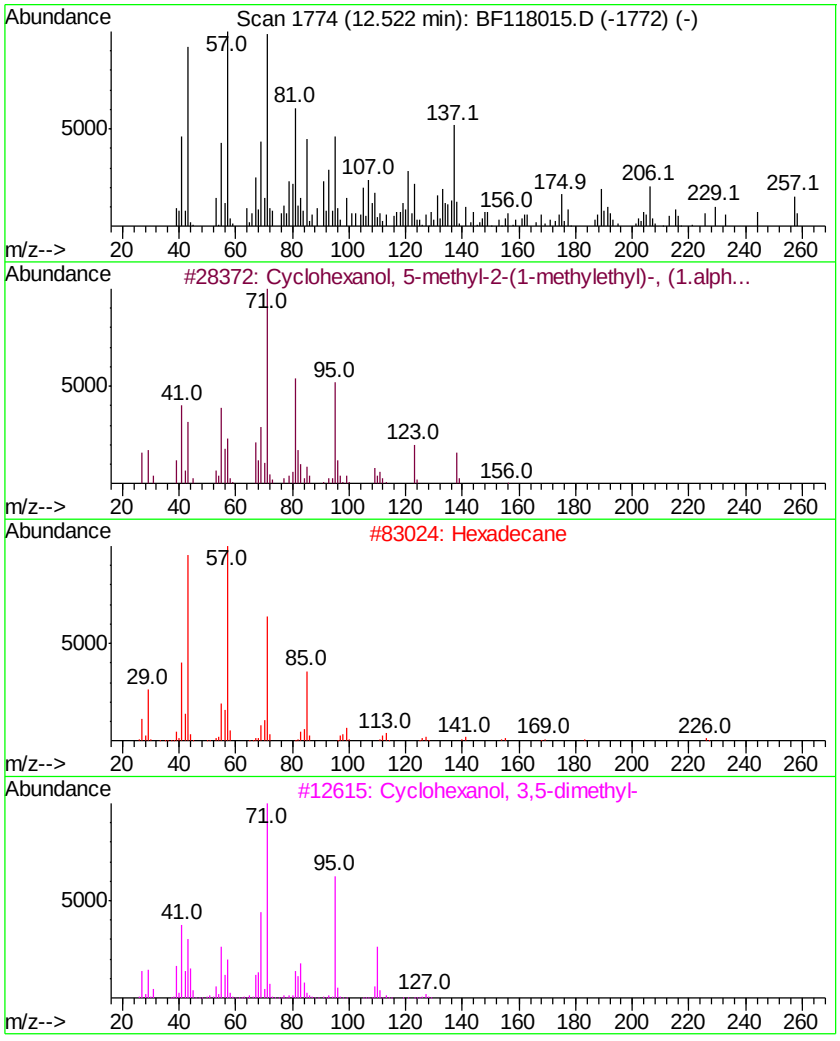
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

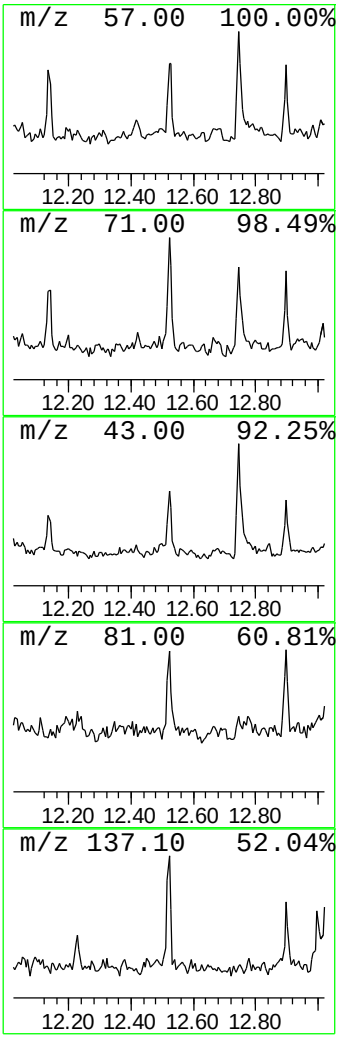
****

Peak Number 9 unknown12.52 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.52	3.00 ng	104699	Phenanthrene-d10	11.44

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanol, 5-methyl-2-(1-meth...	156	C10H20O	000490-99-3	35
2	Hexadecane	226	C16H34	000544-76-3	30
3	Cyclohexanol, 3,5-dimethyl-	128	C8H16O	005441-52-1	30
4	2,6,10,10-Tetramethylbicyclo[7.2...	204	C15H24	136296-37-2	25
5	Sulfurous acid, hexyl pentyl ester	236	C11H24O3S	1000309-14-1	25





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112619\
Data File : BF118015.D
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Instrument : BNA_F
ClientSampleId : B-1-0-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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TIC Integration Parameters: LSCINT.P

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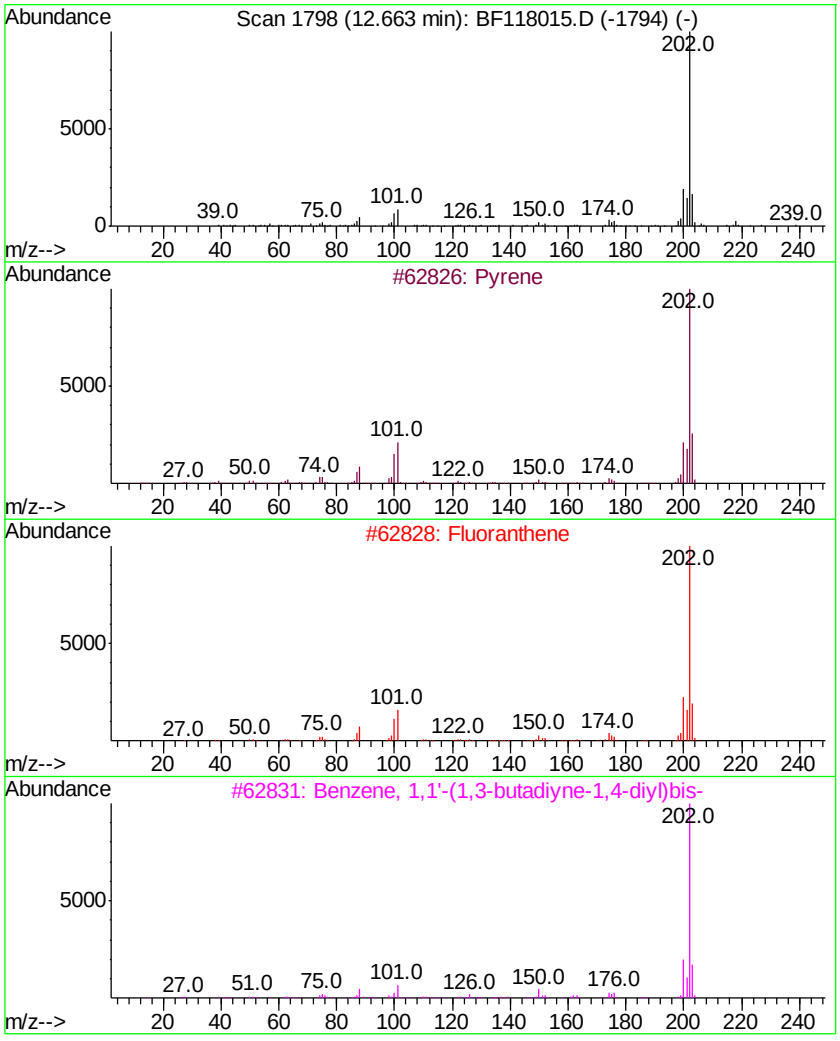
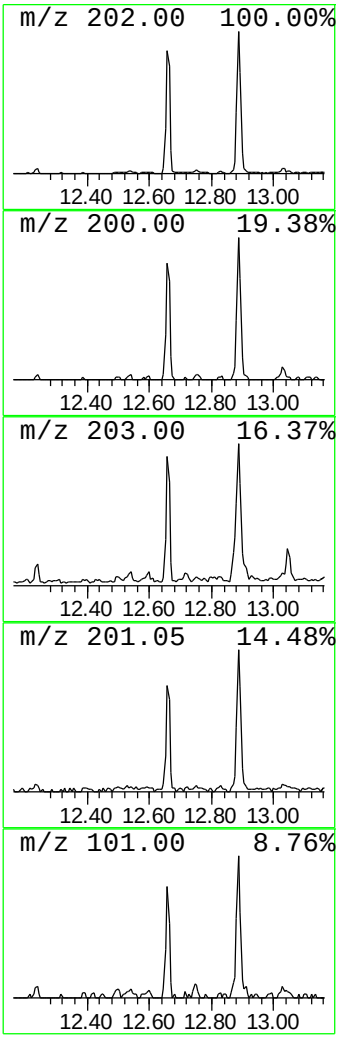
Peak Number 10 Benzene, 1,1'-(1,3-butadiyn... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.66	5.47 ng	191140	Phenanthrene-d10	11.44

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene	202	C16H10	000129-00-0	95
2	Fluoranthene	202	C16H10	000206-44-0	95
3	Benzene, 1,1'-(1,3-butadiyne-1,4...	202	C16H10	000886-66-8	76
4	1,4-Pentadiyn-3-one, 1,5-diphenyl-	230	C17H10O	015814-30-9	64
5	4,4'-Bis(tetrahydrothiopyran)	202	C10H18S2	116196-83-9	60





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112619\
Data File : BF118015.D
Acq On : 26 Nov 2019 18:56
Operator : JU
Sample : K5948-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : B-1-0-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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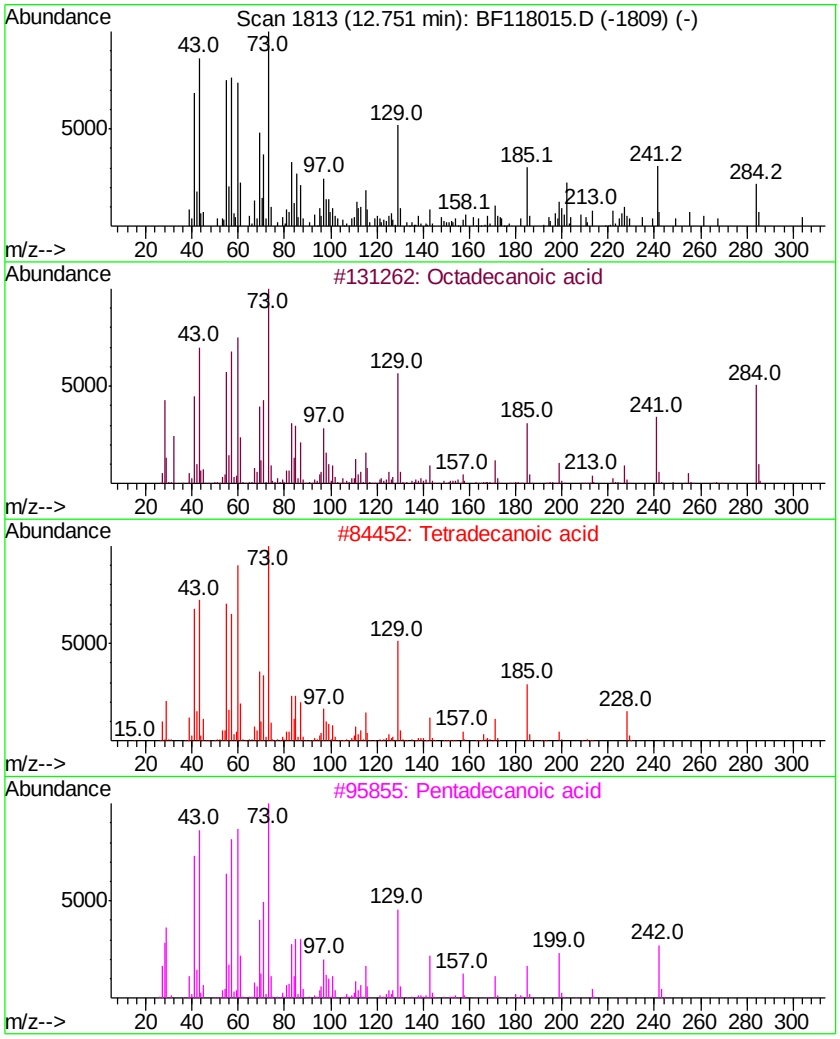
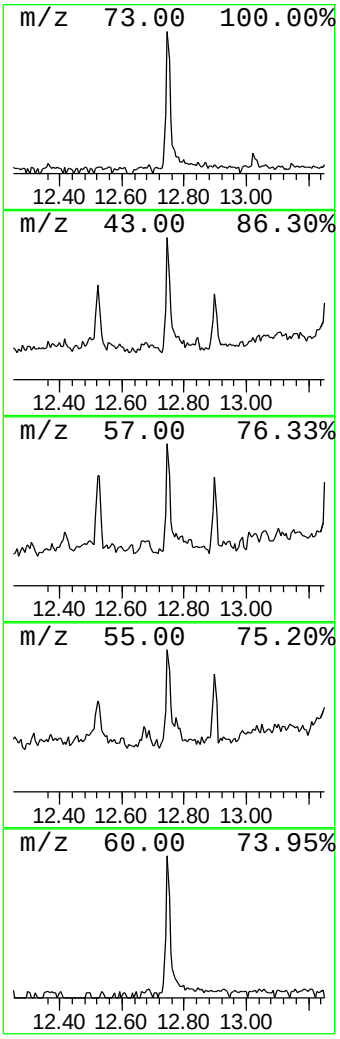
Peak Number 11 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.75	4.07 ng	142130	Phenanthrene-d10	11.44

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	90
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	81
4	Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	80
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	72





Library Search Compound Report

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Operator : JU
Sample : K5948-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : B-1-0-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
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TIC Library : C:\DATABASE\NIST11.L
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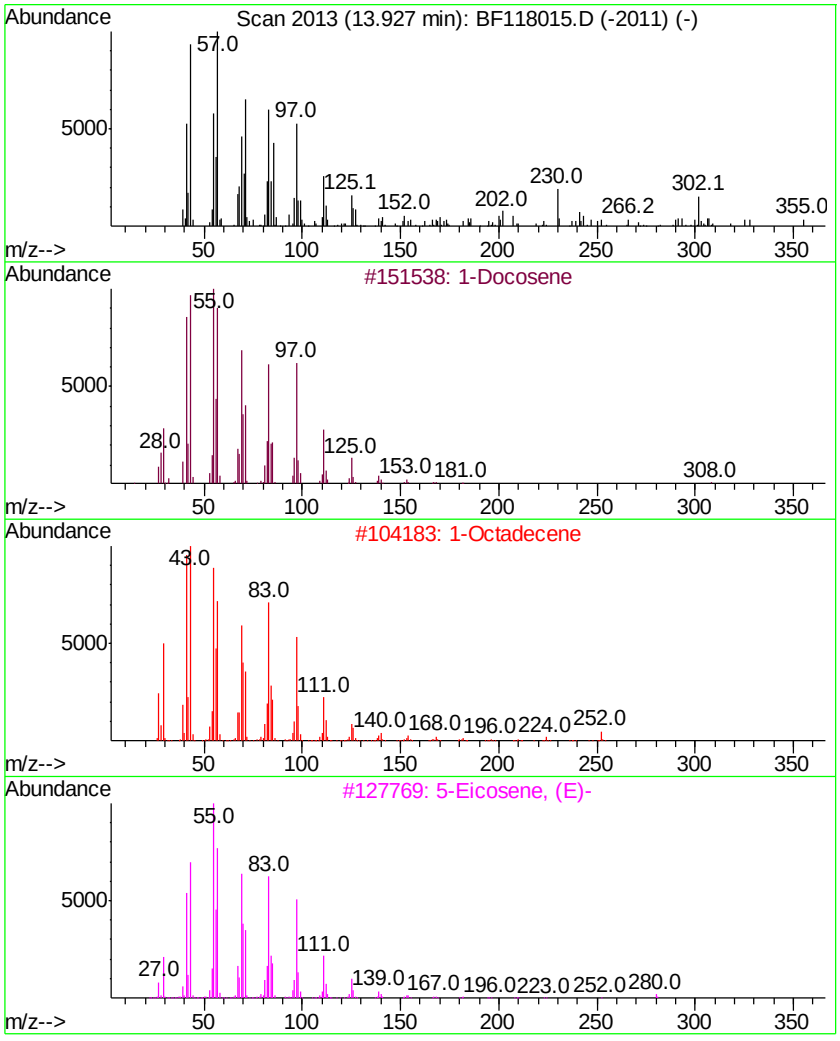
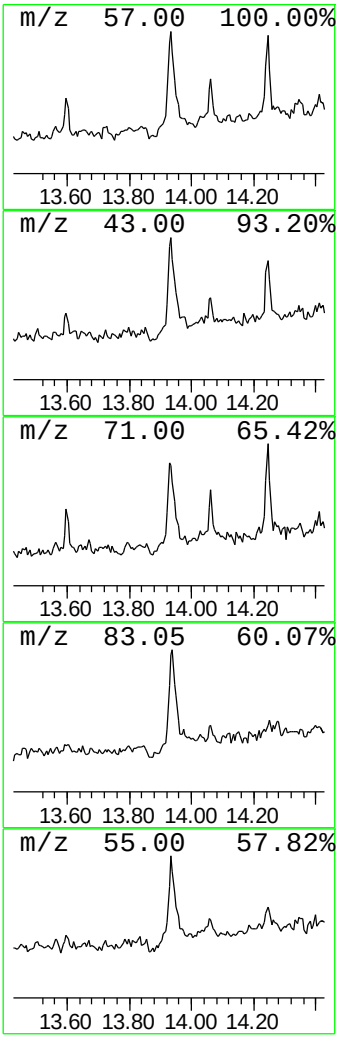
Peak Number 12 1-Docosene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.93	4.78 ng	199013	Chrysene-d12	14.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	94
2	1-Octadecene	252	C18H36	000112-88-9	93
3	5-Eicosene, (E)-	280	C20H40	074685-30-6	92
4	1-Heneicosanol	312	C21H44O	015594-90-8	91
5	Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	89





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Sample : K5948-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : B-1-0-6

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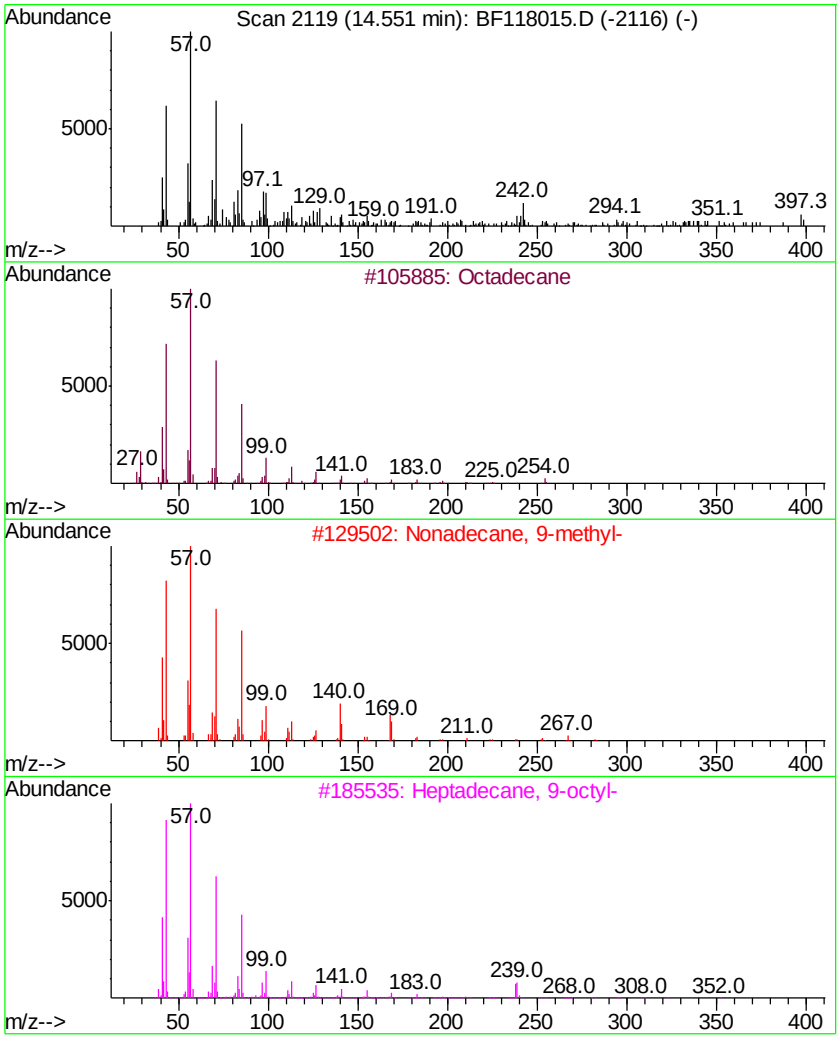
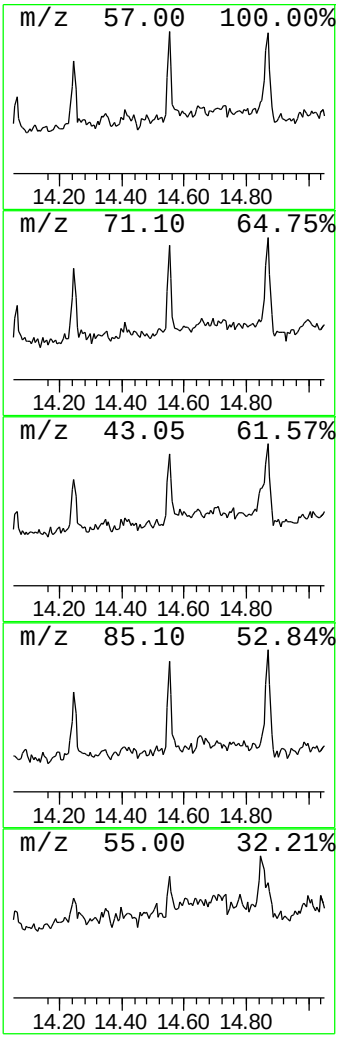
Peak Number 13 Octadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.55	4.42 ng	184102	Chrysene-d12	14.09

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane	254	C18H38	000593-45-3	97
2	Nonadecane, 9-methyl-	282	C20H42	013287-24-6	93
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4	Heptadecane	240	C17H36	000629-78-7	89
5	Hentriacontane	437	C31H64	000630-04-6	81





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ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : B-1-0-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF111319.M
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TIC Integration Parameters: LSCINT.P

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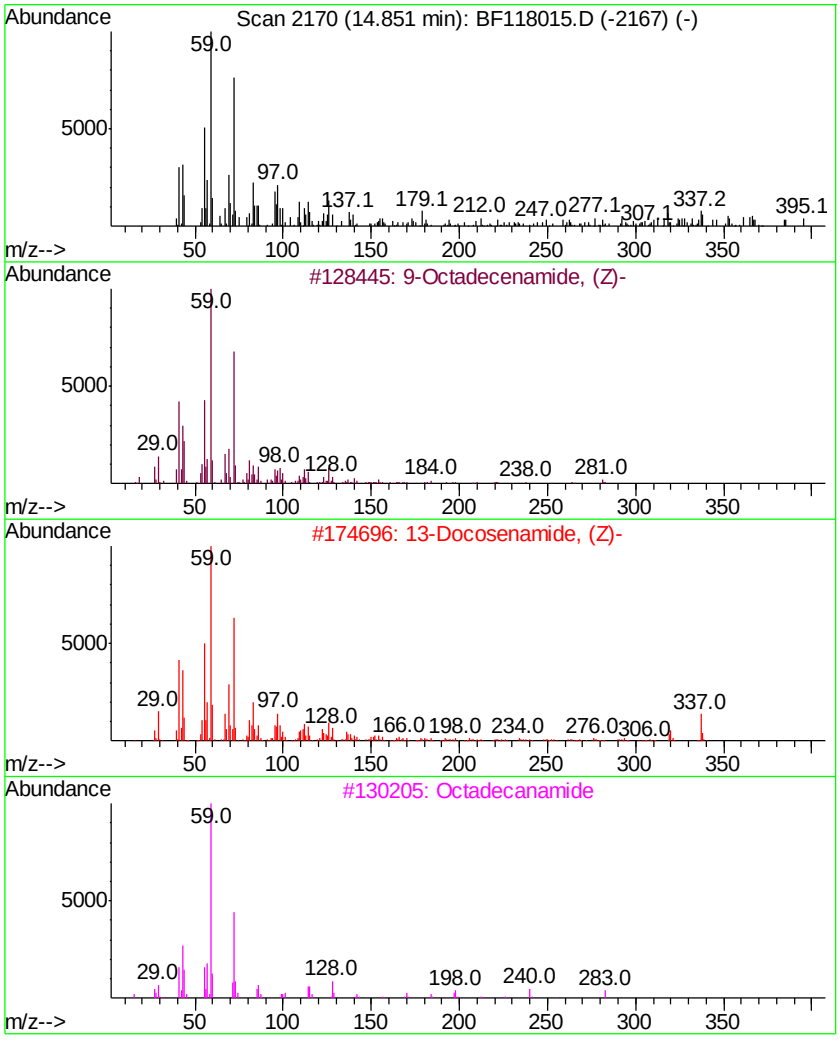
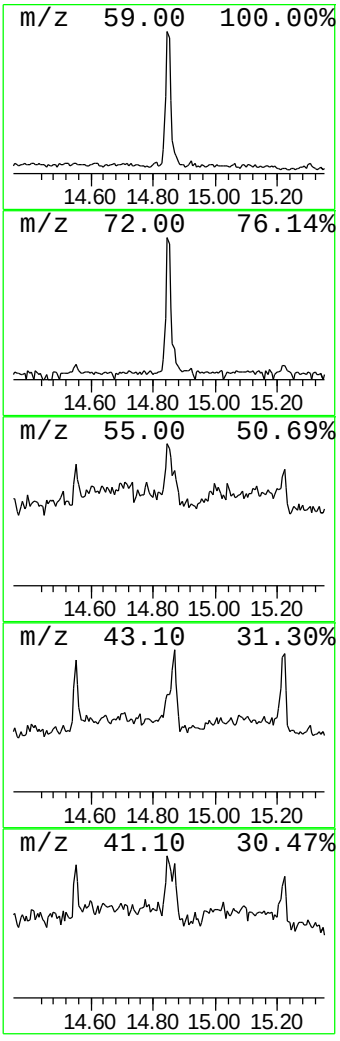
Peak Number 14 9-Octadecenamide, (Z)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.85	6.29 ng	259310	Perylene-d12	15.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	93
2	13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	81
3	Octadecanamide	283	C18H37NO	000124-26-5	43
4	Tetradecanamide	227	C14H29NO	000638-58-4	35
5	Dodecanamide	199	C12H25NO	001120-16-7	27





Tentatively Identified Compound (LSC) summary

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ClientSampleId : B-1-0-6

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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...	2.19	20.9	ng	680750	1	6.89	650916	20.0
2-Pentanone, 4-hy...	5.11	16.1	ng	522291	1	6.89	650916	20.0
unknown6.66	6.66	76.6	ng	2494140	1	6.89	650916	20.0
Hexadecane, 2,6,1...	10.86	4.3	ng	149065	4	11.44	698899	20.0
n-Hexadecanoic acid	11.96	11.8	ng	413447	4	11.44	698899	20.0
unknown12.00	12.00	2.5	ng	86902	4	11.44	698899	20.0
unknown12.52	12.52	3.0	ng	104699	4	11.44	698899	20.0
Benzene, 1,1'-(1,...	12.66	5.5	ng	191140	4	11.44	698899	20.0
Octadecanoic acid	12.75	4.1	ng	142130	4	11.44	698899	20.0
1-Docosene	13.93	4.8	ng	199013	5	14.09	833123	20.0
Octadecane	14.55	4.4	ng	184102	5	14.09	833123	20.0
9-Octadecenamide,...	14.85	6.3	ng	259310	6	15.60	825143	20.0