LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF123019\
Data File : BF118390.D
Acq On : 30 Dec 2019 19:33
Operator : JU
Sample : K6447-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RBR31-48

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-BF122419.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	5.057	501	505	519	rVB	341675	458651	11.96%	1.209%
2	5.475	572	576	599	rBV	2372903	2517950	65.65%	6.635%
3	6.492	744	749	767	rBV	2479477	2697788	70.34%	7.109%
4	6.622	767	771	775	rBV	3065505	2902313	75.67%	7.648%
5	6.851	806	810	821	rVB	913290	735133	19.17%	1.937%
6	7.010	833	837	840	rBV	2675144	2309733	60.22%	6.086%
7	7.416	902	906	929	rBV	1620001	1742316	45.43%	4.591%
8	8.139	1025	1029	1047	rBV	977595	975395	25.43%	2.570%
9	9.216	1208	1212	1223	rBV	4259480	3708389	96.68%	9.772%
10	9.551	1267	1269	1274	rVB2	127303	98454	2.57%	0.259%
11	9.616	1277	1280	1285	rVB	175602	183566	4.79%	0.484%
12	9.869	1319	1323	1325	rBV	79104	63441	1.65%	0.167%
13	9.898	1325	1328	1332	rVV	1263567	1163076	30.32%	3.065%
14	9.933	1332	1334	1341	rVB	60601	63506	1.66%	0.167%
15	10.110	1361	1364	1368	rBV2	36749	44253	1.15%	0.117%
16	10.698	1459	1464	1479	rVB	2378581	2509889	65.44%	6.614%
17	11.051	1521	1524	1533	rBV	179036	199518	5.20%	0.526%
18	11.392	1578	1582	1585	rBV	1369105	1241760	32.37%	3.272%
19	11.416	1585	1586	1591	rVB	261739	184979	4.82%	0.487%
20	11.857	1656	1661	1664	rBV3	35126	64179	1.67%	0.169%
21	11.927	1665	1673	1685	rVV	1944862	2383128	62.13%	6.280%
22	12.545	1774	1778	1781	rBV3	35021	44560	1.16%	0.117%
23	12.616	1786	1790	1791	rBV	472011	525883	13.71%	1.386%
24	12.633	1791	1793	1803	rVV	975343	1135878	29.61%	2.993%
25	12.710	1803	1806	1814	rVB	620932	646689	16.86%	1.704%
26	12.845	1823	1829	1835	rVB	297298	324129	8.45%	0.854%
27	12.986	1848	1853	1856	rBV	4489758	3835561	100.00%	10.107%
28	13.745	1979	1982	1986	rBV	257412	224477	5.85%	0.592%
29	13.880	2001	2005	2016	rVB2	477389	698227	18.20%	1.840%
30	14.051	2029	2034	2037	rBV	954504	1000256	26.08%	2.636%
31	14.074	2037	2038	2042	rVB	108138	70017	1.83%	0.184%
32	14.198	2055	2059	2065	rBV	84287	105498	2.75%	0.278%
33	14.504	2108	2111	2113	rBV	152181	151921	3.96%	0.400%
34	14.809	2159	2163	2166	rBV	116942	123260	3.21%	0.325%

LSC Area Percent Report

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Acq On : 30 Dec 2019 19:33
Operator : JU
Sample : K6447-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RBR31-48

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-BF122419.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	14.886	2173	2176	2180	rBV	53951	53962	1.41%	0.142%
36	14.962	2186	2189	2193	rVB4	51328	57734	1.51%	0.152%
37	15.104	2210	2213	2217	rBV	90430	142168	3.71%	0.375%
38	15.157	2218	2222	2225	rVV	205628	238106	6.21%	0.627%
39	15.198	2225	2229	2238	rVV7	64533	147929	3.86%	0.390%
40	15.415	2263	2266	2270	rBV	65871	75900	1.98%	0.200%
41	15.480	2273	2277	2281	rVV	93351	127413	3.32%	0.336%
42	15.545	2283	2288	2299	rVB	774479	1046995	27.30%	2.759%
43	15.745	2319	2322	2327	rVB3	58517	70851	1.85%	0.187%
44	15.986	2358	2363	2369	rVB	176094	266932	6.96%	0.703%
45	16.051	2369	2374	2386	rBV6	96639	253742	6.62%	0.669%
46	16.492	2446	2449	2453	rBV4	37487	59121	1.54%	0.156%
47	16.792	2494	2500	2507	rBV7	49132	106638	2.78%	0.281%
48	17.068	2542	2547	2549	rBV2	26986	49779	1.30%	0.131%
49	17.633	2638	2643	2654	rVB2	43938	119176	3.11%	0.314%

Sum of corrected areas: 37950219

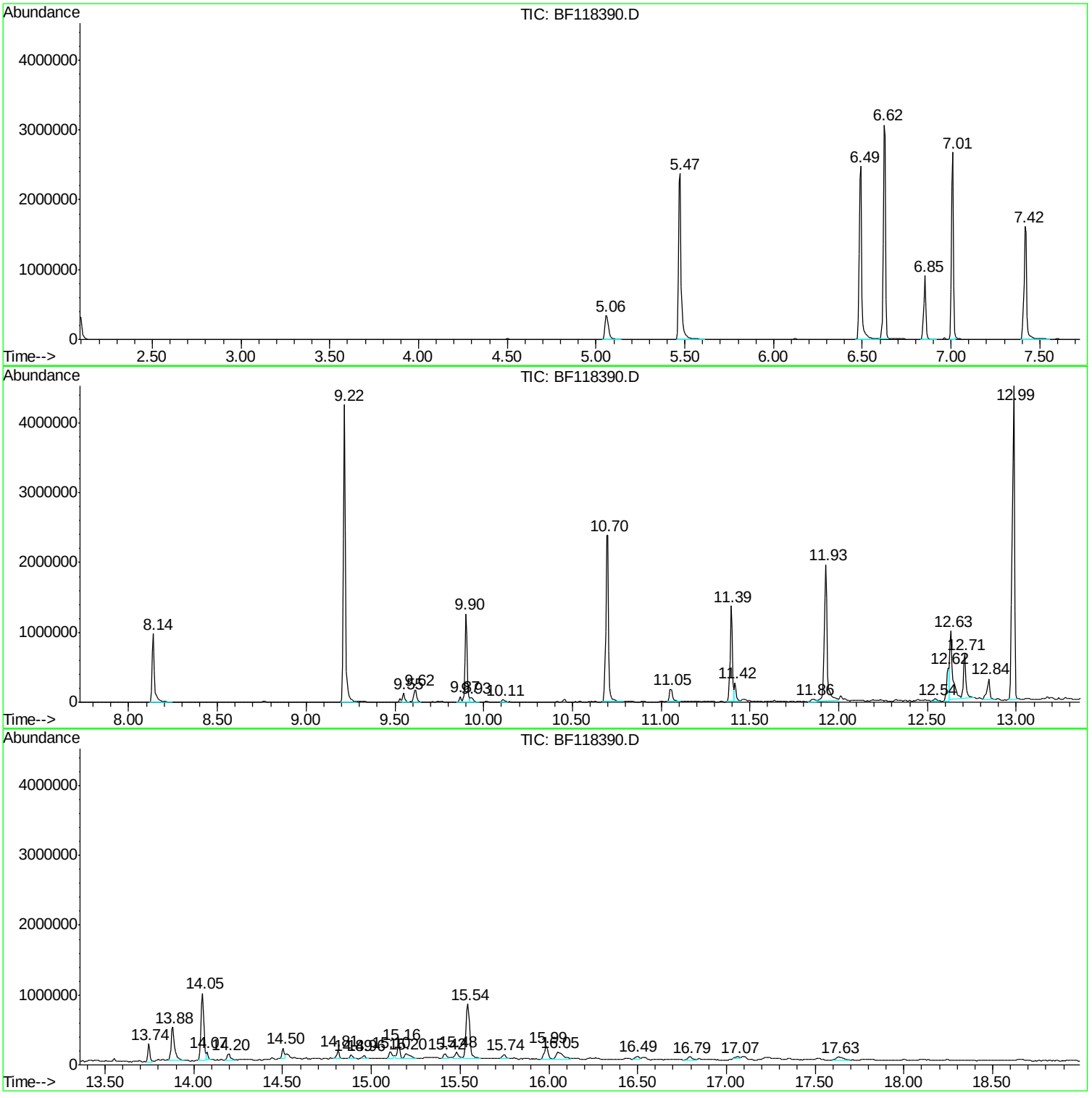
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF123019\
Data File : BF118390.D
Acq On : 30 Dec 2019 19:33
Operator : JU
Sample : K6447-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RBR31-48

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF122419.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\BF123019\
Data File : BF118390.D
Acq On : 30 Dec 2019 19:33
Operator : JU
Sample : K6447-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RBR31-48

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF122419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

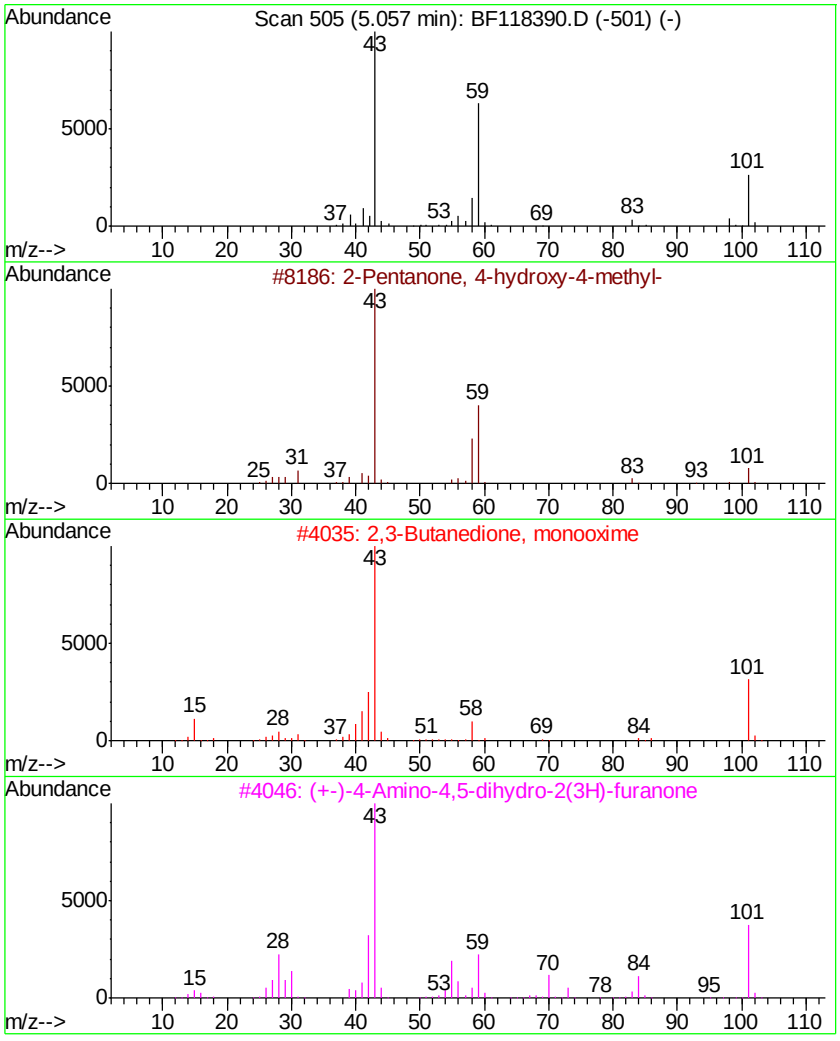
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

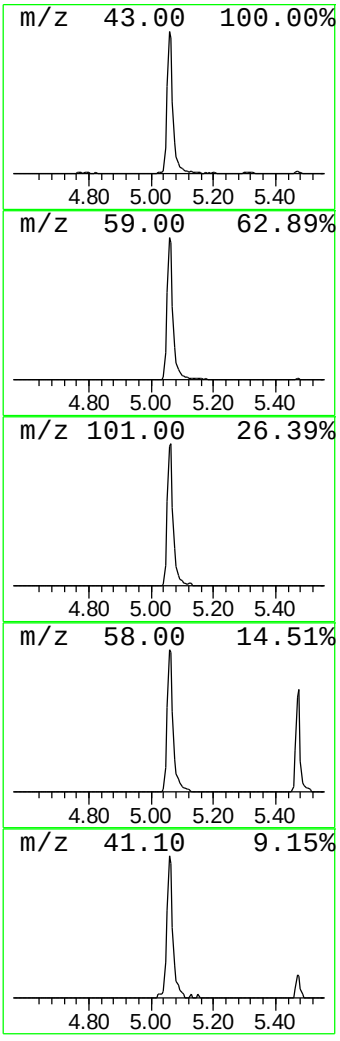
****

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.06	12.48 ng	458651	1,4-Dichlorobenzene-d4	6.85

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	72
2	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	25
3	(+-)-4-Amino-4,5-dihydro-2(3H)-f...	101	C4H7NO2	016504-58-8	9
4	5-Hexen-2-one	98	C6H10O	000109-49-9	9
5	4-Methoxy-1-pentene	100	C6H12O	098386-09-5	9





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF123019\
Data File : BF118390.D
Acq On : 30 Dec 2019 19:33
Operator : JU
Sample : K6447-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RBR31-48

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF122419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

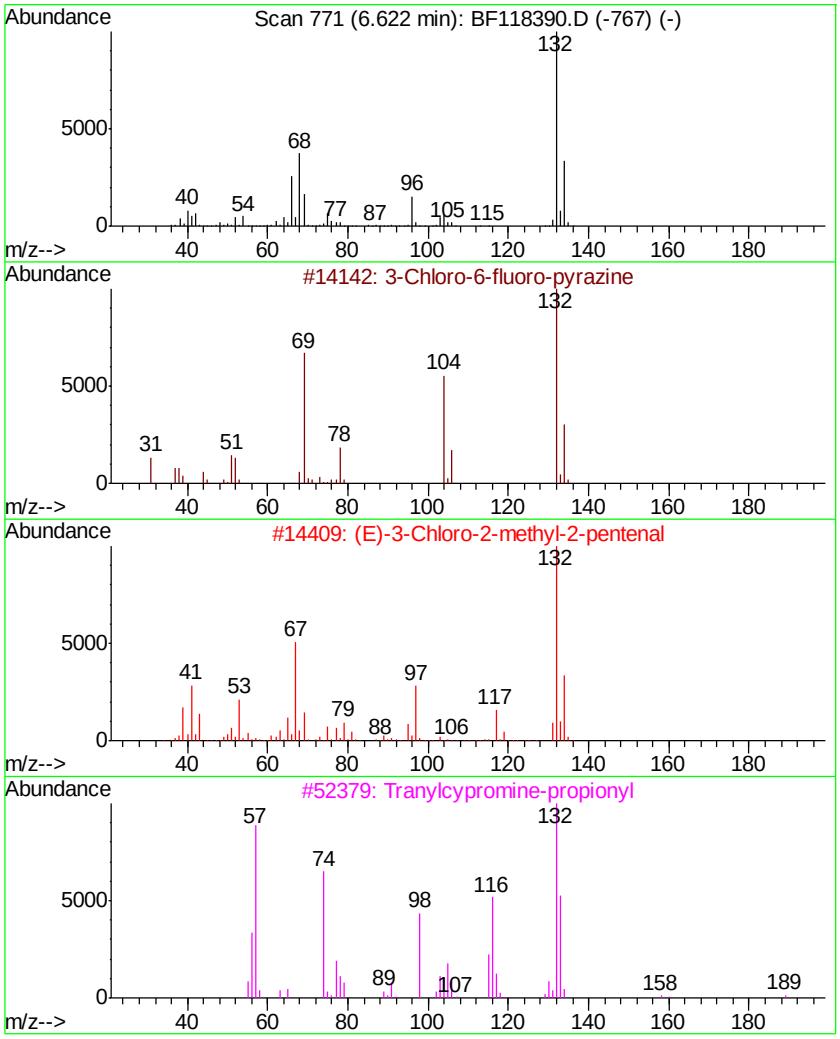
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

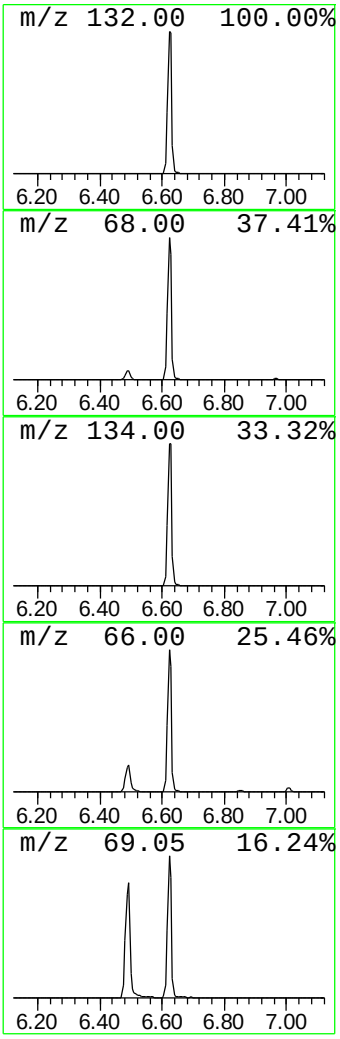
****

Peak Number 2 unknown6.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	78.96 ng	2902310	1,4-Dichlorobenzene-d4	6.85

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	25
3	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4	1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5	Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF123019\
Data File : BF118390.D
Acq On : 30 Dec 2019 19:33
Operator : JU
Sample : K6447-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RBR31-48

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF122419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

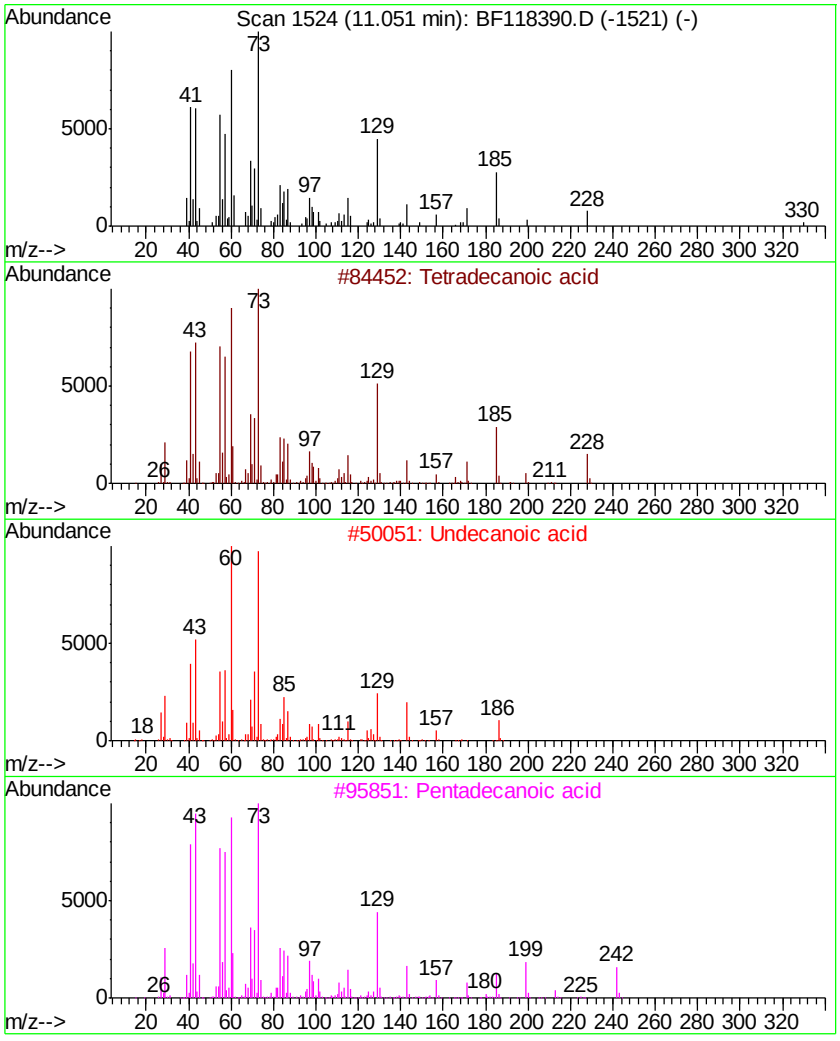
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

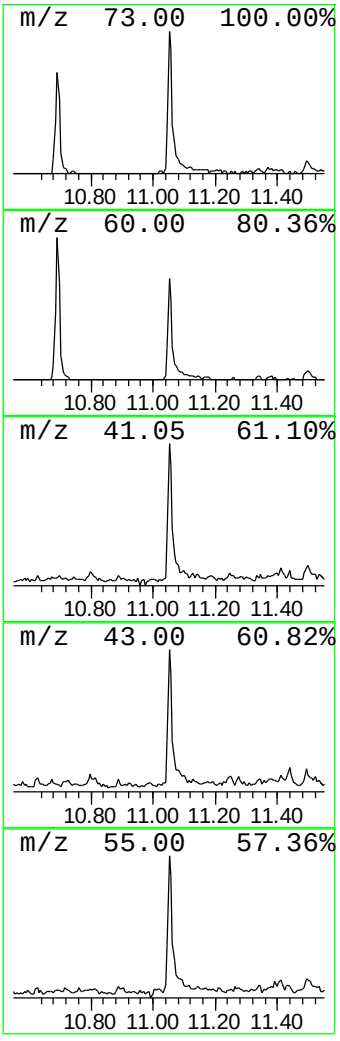
****

Peak Number 4 Tetradecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.05	3.21 ng	199518	Phenanthrene-d10	11.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetradecanoic acid	228	C14H28O2	000544-63-8	98
2	Undecanoic acid	186	C11H22O2	000112-37-8	81
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	76
4	Decanoic acid, silver(1+) salt	278	C10H19AgO2	013126-67-5	72
5	Dodecanoic acid	200	C12H24O2	000143-07-7	72





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF123019\
Data File : BF118390.D
Acq On : 30 Dec 2019 19:33
Operator : JU
Sample : K6447-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RBR31-48

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF122419.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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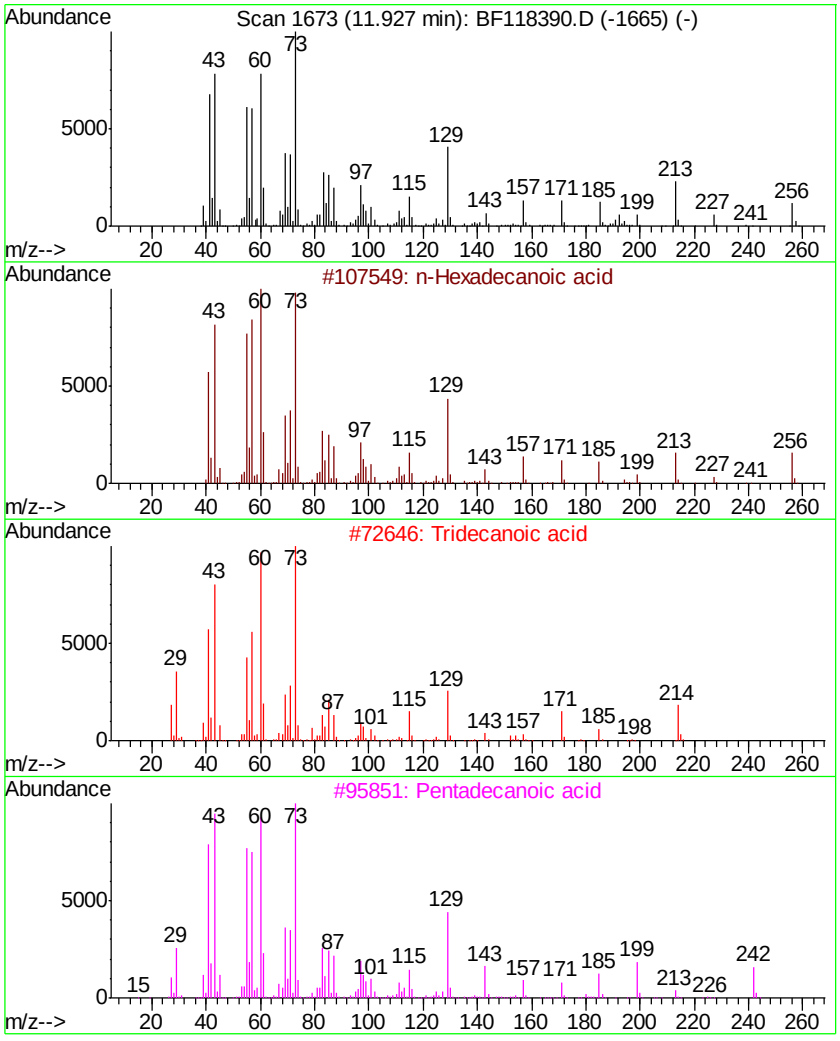
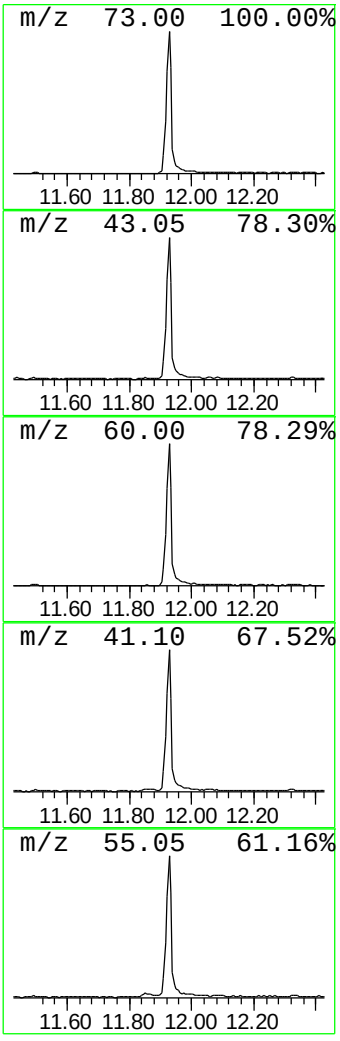
Peak Number 5 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	38.38 ng	2383130	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	91
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	87
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	72
5	n-Decanoic acid	172	C10H20O2	000334-48-5	58





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF123019\
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Acq On : 30 Dec 2019 19:33
Operator : JU
Sample : K6447-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RBR31-48

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF122419.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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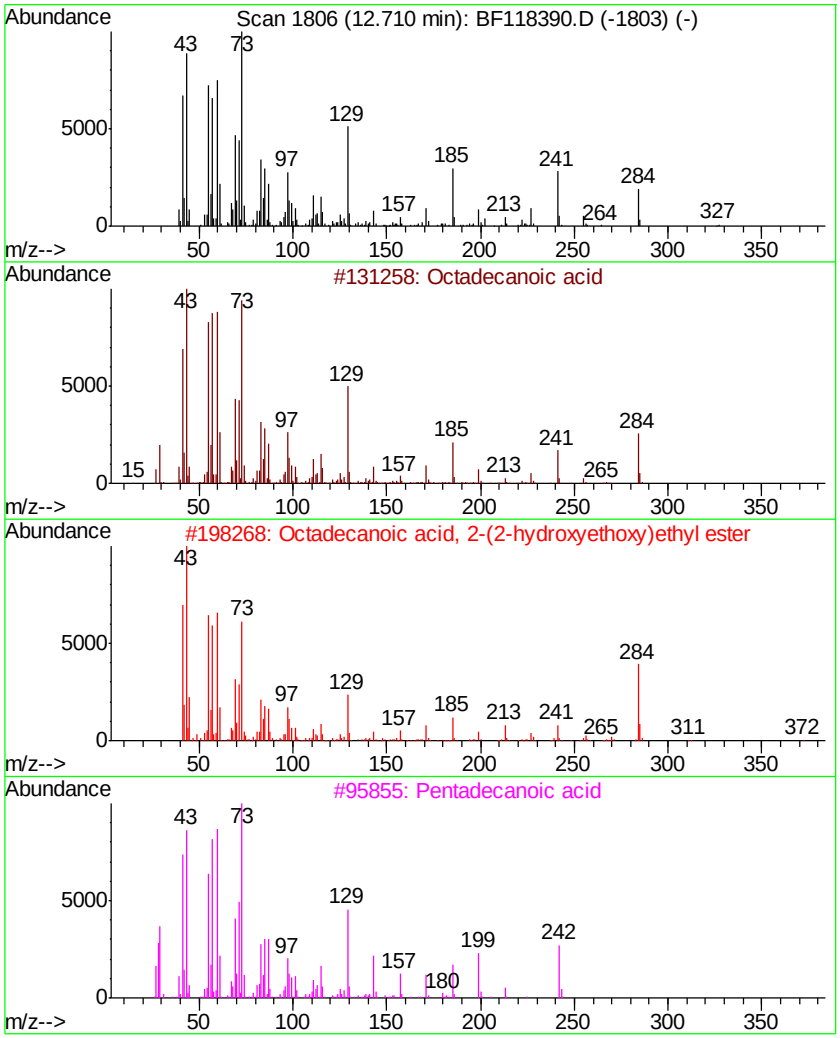
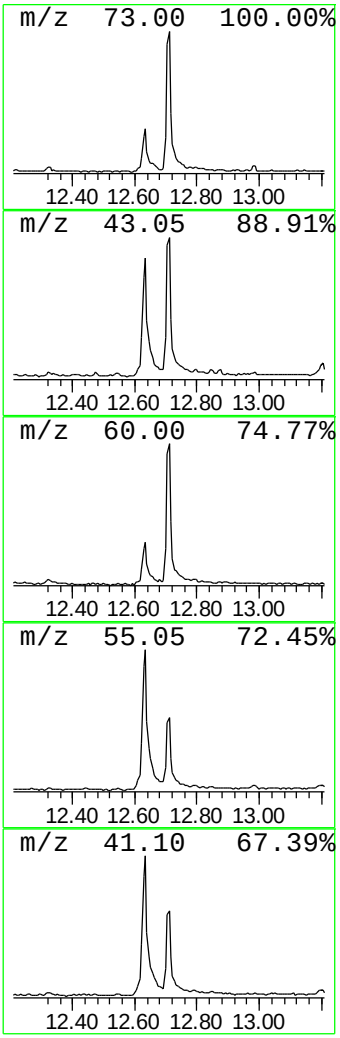
Peak Number 6 Octadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.71	10.42 ng	646689	Phenanthrene-d10	11.39

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	91
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	89
4	n-Decanoic acid	172	C10H20O2	000334-48-5	64
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	62





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Data File : BF118390.D
Acq On : 30 Dec 2019 19:33
Operator : JU
Sample : K6447-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RBR31-48

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF122419.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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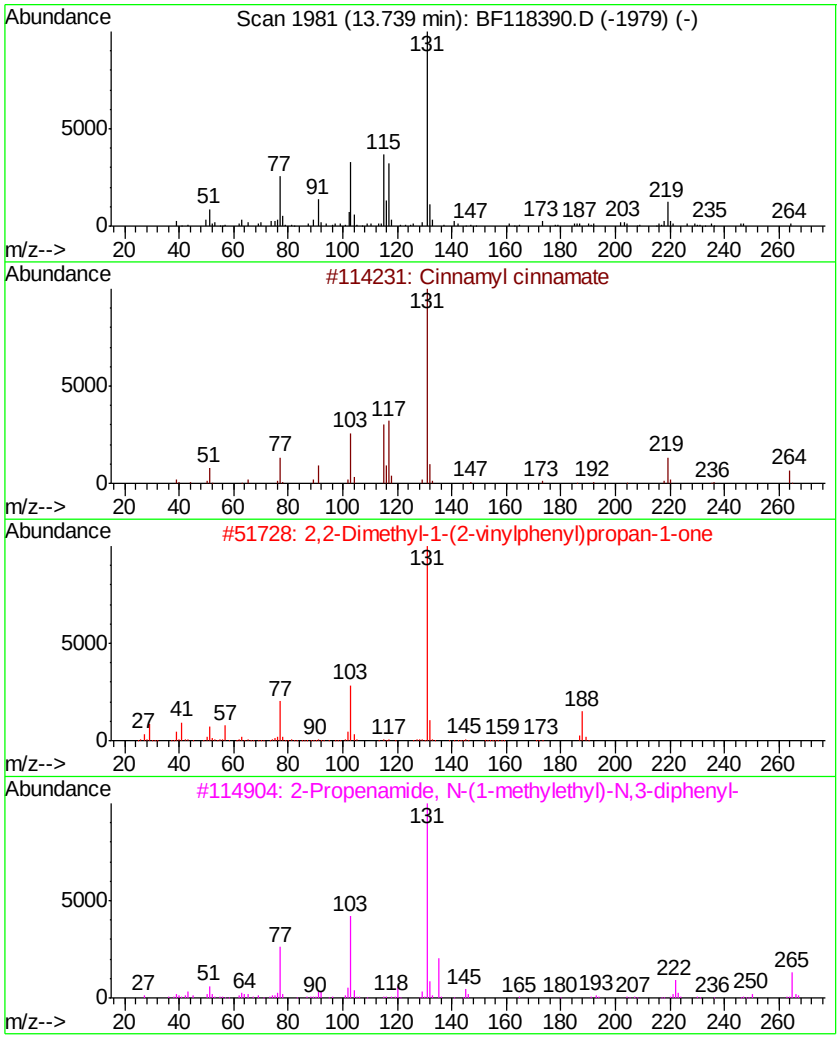
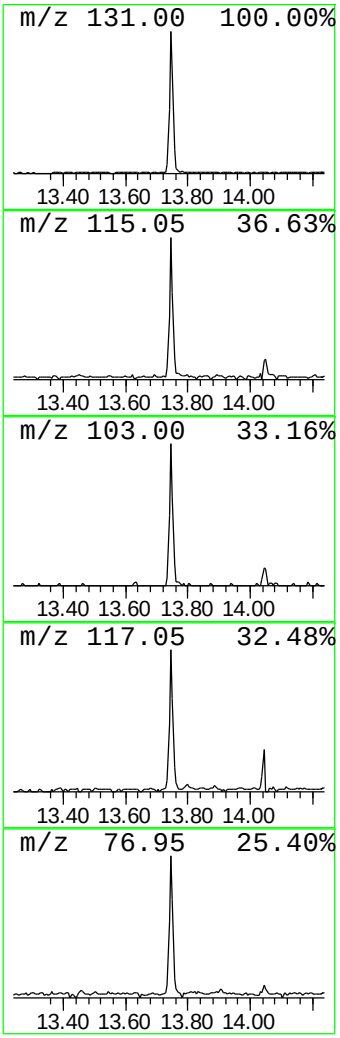
Peak Number 7 Cinnamyl cinnamate Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.74	4.49 ng	224477	Chrysene-d12	14.05

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cinnamyl cinnamate	264	C18H16O2	000122-69-0	87
2	2,2-Dimethyl-1-(2-vinylphenyl)pr...	188	C13H16O	1000210-99-9	52
3	2-Propenamide, N-(1-methylethyl)...	265	C18H19NO	055044-37-6	50
4	2-Propenamide, N-(1,1-dimethylet...	203	C13H17NO	003579-53-1	50
5	2-Propenamide, N,N-diethyl-3-phe...	203	C13H17NO	003680-04-4	50





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Data File : BF118390.D
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Operator : JU
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Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RBR31-48

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF122419.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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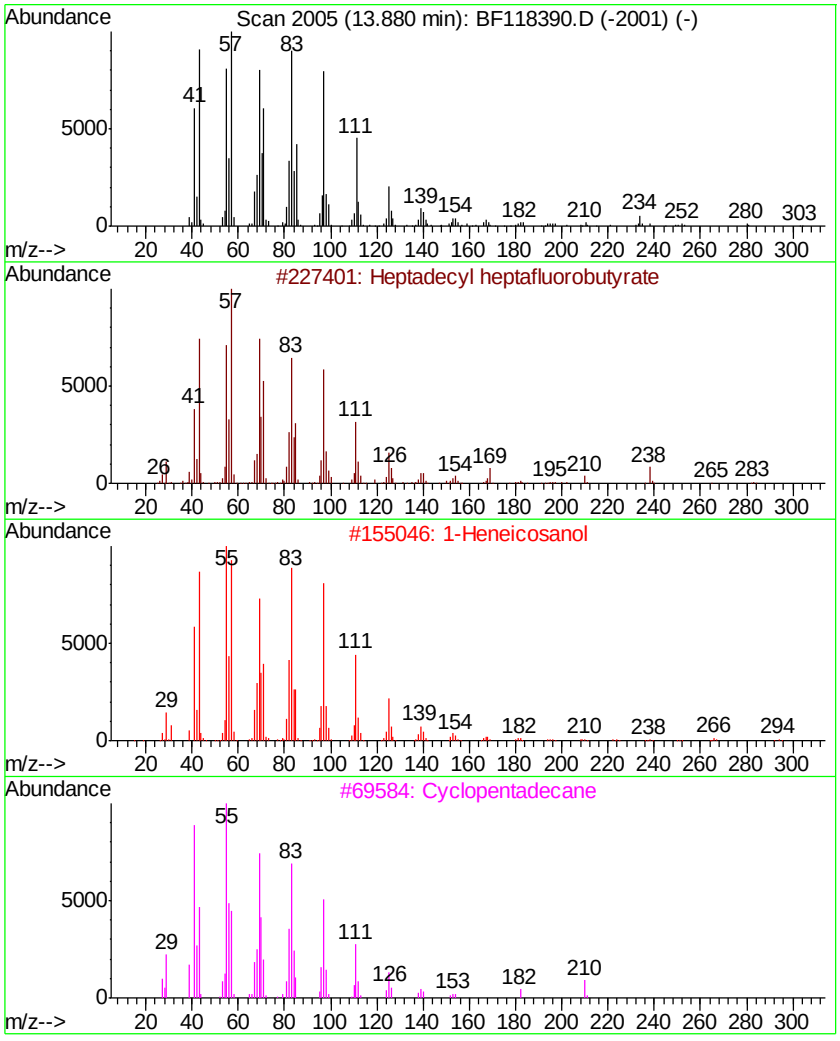
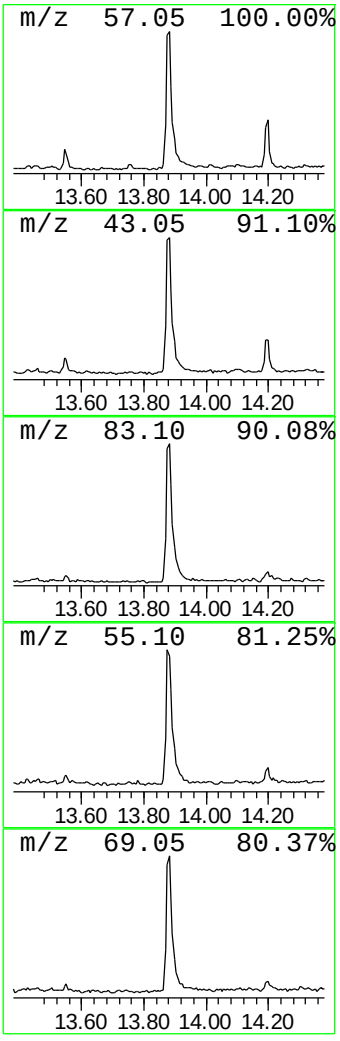
Peak Number 8 Heptadecyl heptafluorobutyrate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.88	13.96 ng	698227	Chrysene-d12	14.05

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
2	1-Heneicosanol	312	C21H44O	015594-90-8	93
3	Cyclopentadecane	210	C15H30	000295-48-7	93
4	n-Tetracosanol-1	354	C24H50O	000506-51-4	93
5	Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	92





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF123019\
Data File : BF118390.D
Acq On : 30 Dec 2019 19:33
Operator : JU
Sample : K6447-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RBR31-48

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF122419.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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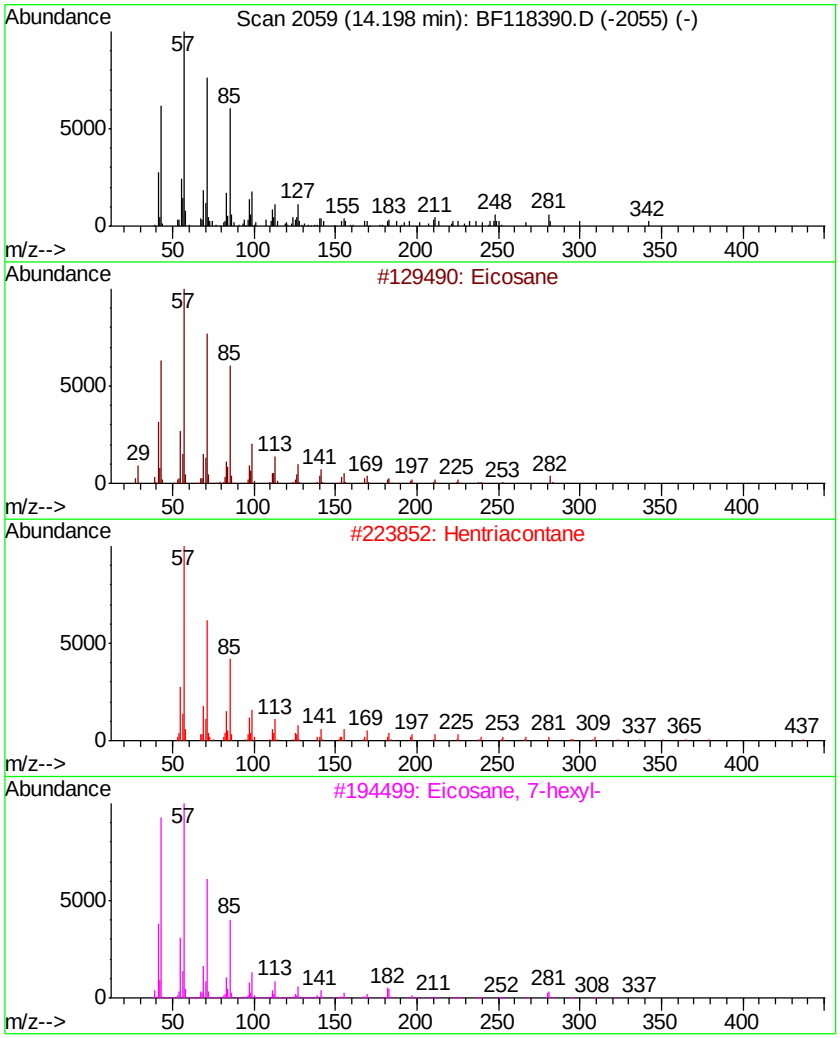
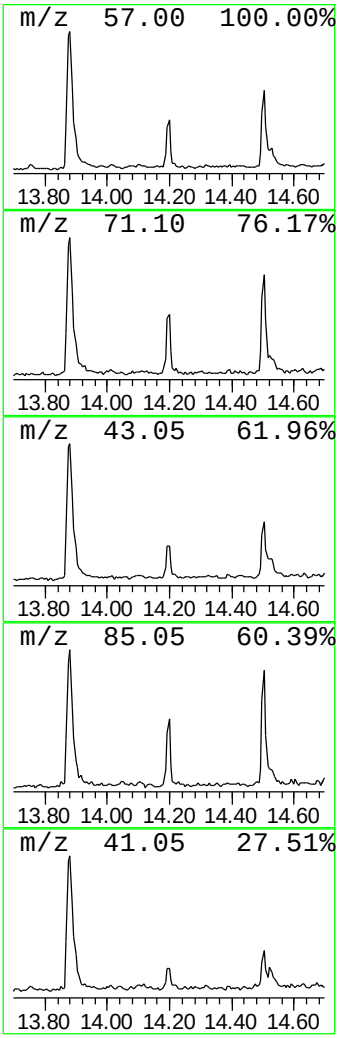
Peak Number 9 Eicosane Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.20	2.11 ng	105498	Chrysene-d12	14.05

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	93
2	Hentriacontane	437	C31H64	000630-04-6	91
3	Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
4	Hexacosane	366	C26H54	000630-01-3	90
5	Heneicosane	296	C21H44	000629-94-7	87





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Acq On : 30 Dec 2019 19:33
Operator : JU
Sample : K6447-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RBR31-48

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF122419.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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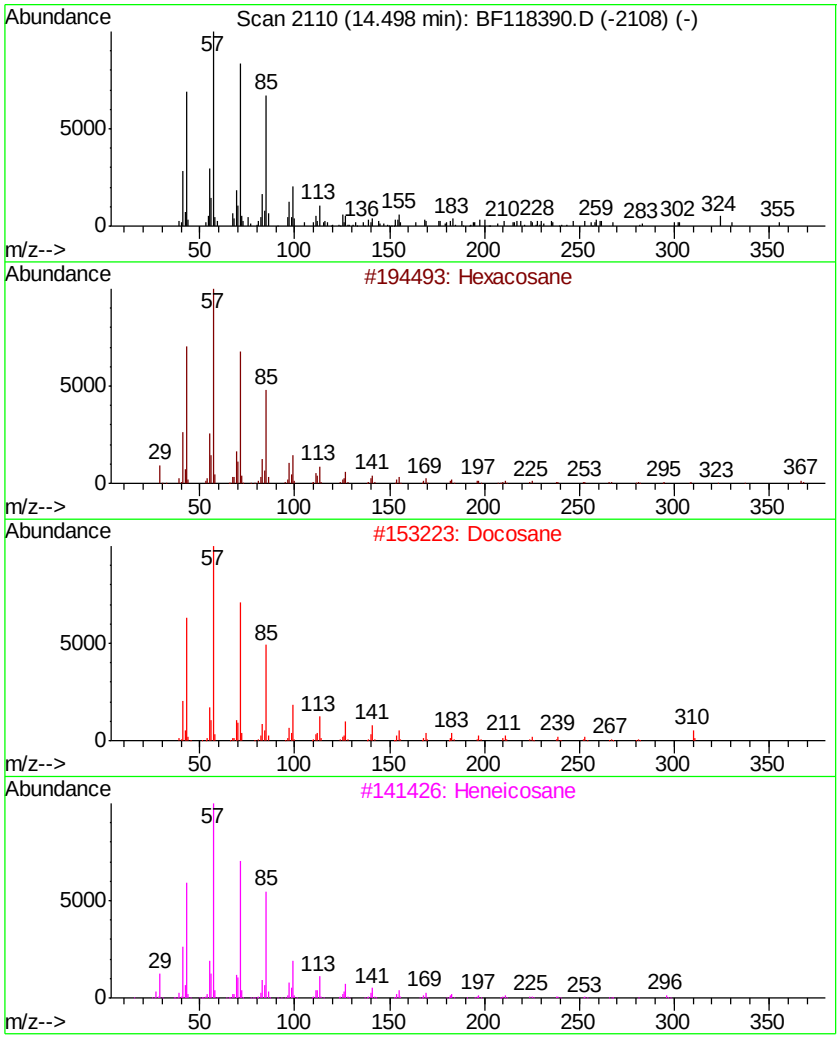
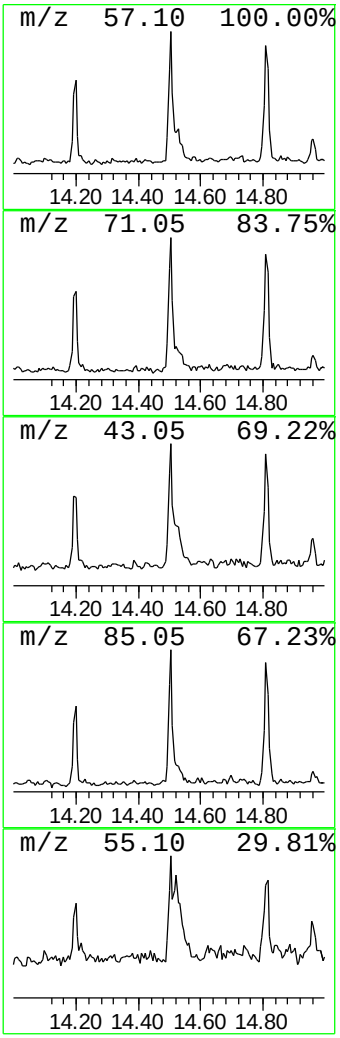
Peak Number 10 Hexacosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.50	3.04 ng	151921	Chrysene-d12	14.05

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	90
2	Docosane	310	C22H46	000629-97-0	90
3	Heneicosane	296	C21H44	000629-94-7	90
4	Eicosane	282	C20H42	000112-95-8	90
5	Triacontane	422	C30H62	000638-68-6	90





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Instrument : BNA_F
ClientSampleId : RBR31-48

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

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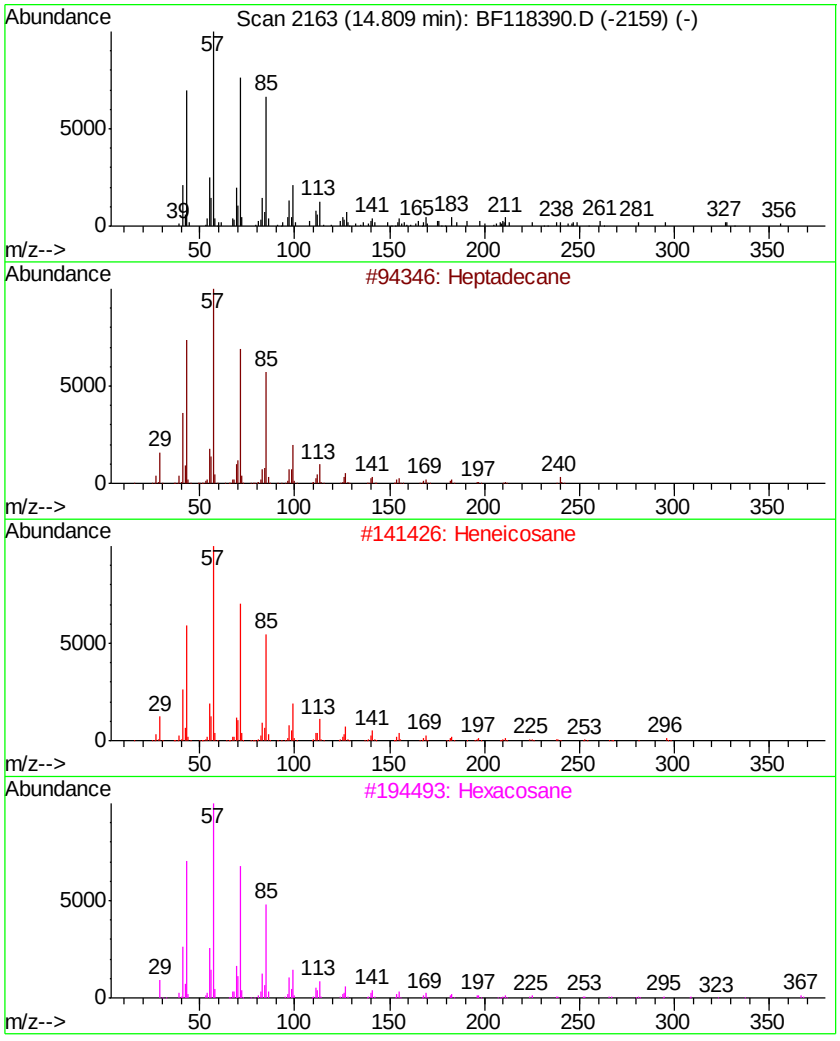
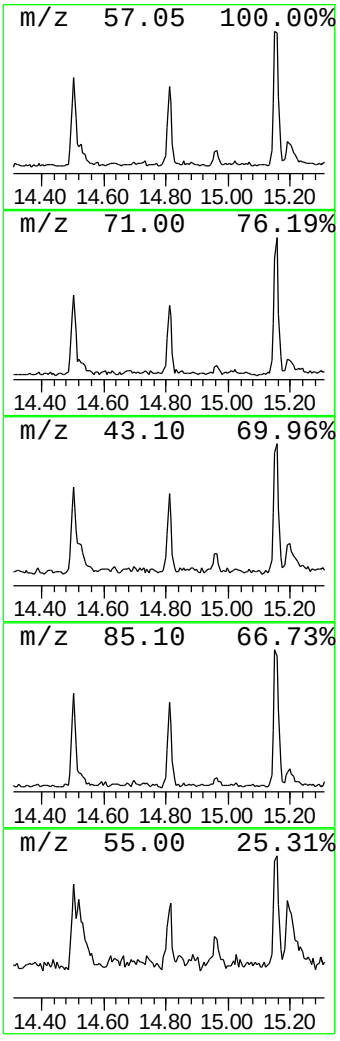
Peak Number 11 Heptadecane Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.81	2.35 ng	123260	Perylene-d12	15.54

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	93
2	Heneicosane	296	C21H44	000629-94-7	90
3	Hexacosane	366	C26H54	000630-01-3	87
4	Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	87
5	Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	81





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Operator : JU
Sample : K6447-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RBR31-48

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TIC Library : C:\DATABASE\NIST11.L
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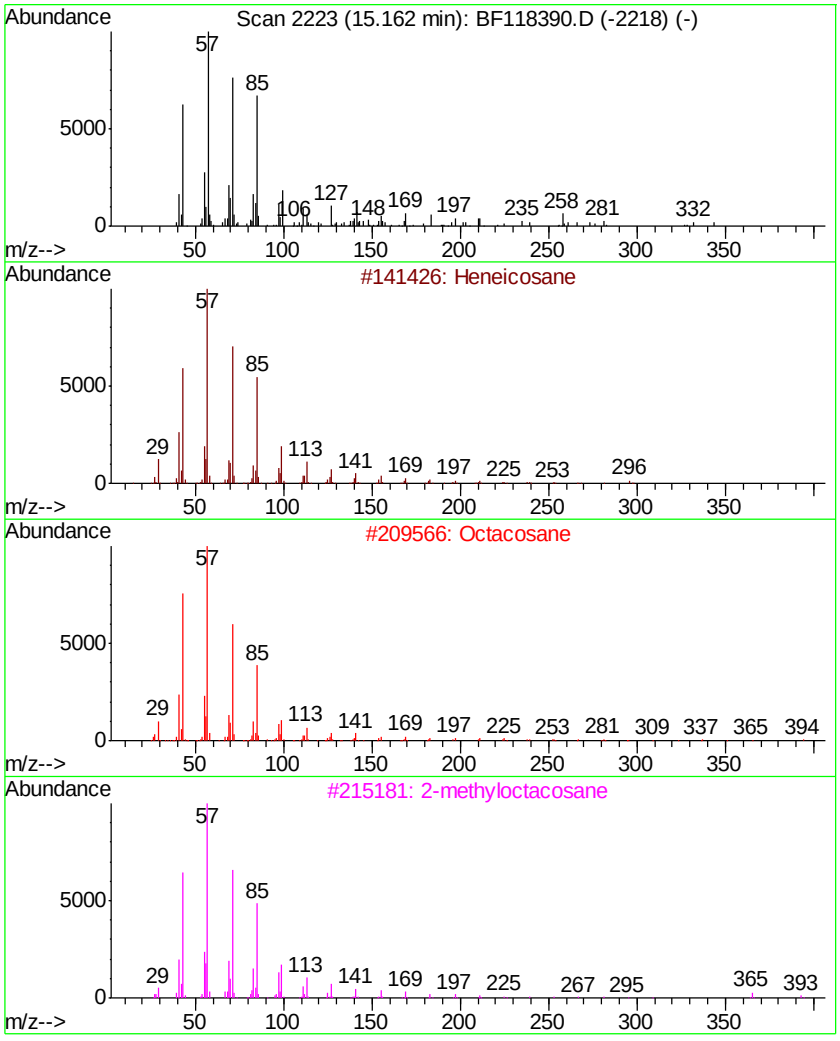
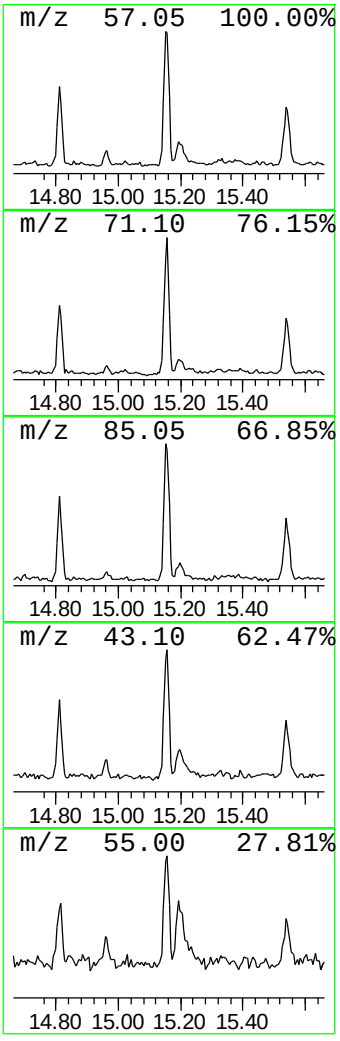
Peak Number 12 Heneicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.16	4.55 ng	238106	Perylene-d12	15.54

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	Octacosane	394	C28H58	000630-02-4	91
3	2-methyloctacosane	408	C29H60	1000376-72-8	91
4	Heptacosane	380	C27H56	000593-49-7	91
5	Docosane, 11-decyl-	451	C32H66	055401-55-3	91





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Operator : JU
Sample : K6447-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RBR31-48

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF122419.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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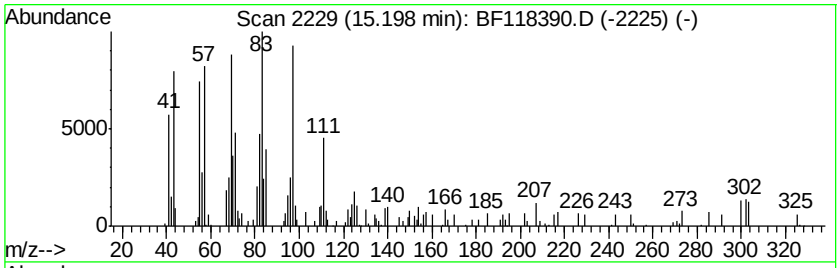
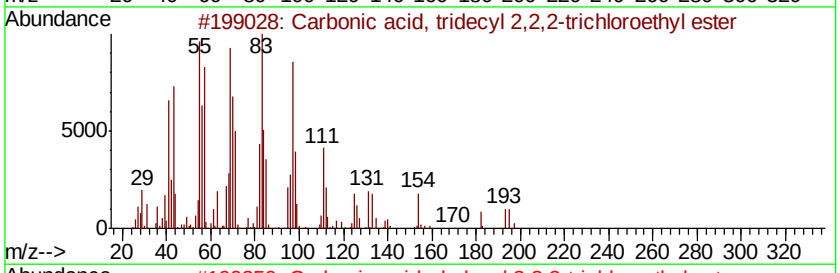
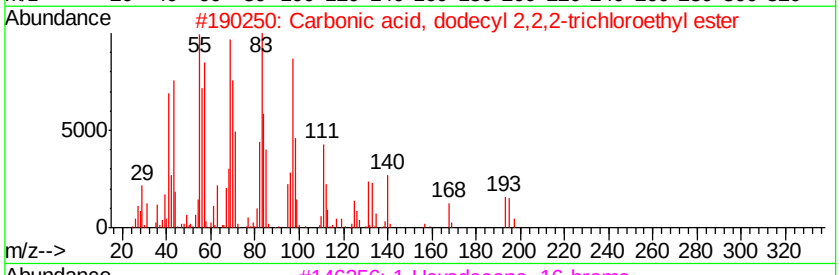
Peak Number 13 Carbonic acid, tridecyl 2,2... Concentration Rank 13

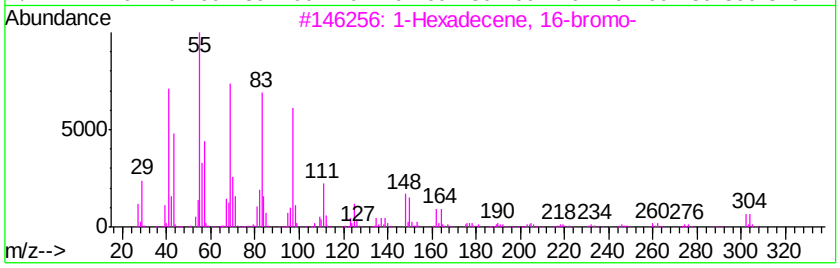
R.T.	EstConc	Area	Relative to ISTD	R.T.
15.20	2.83 ng	147929	Perylene-d12	15.54

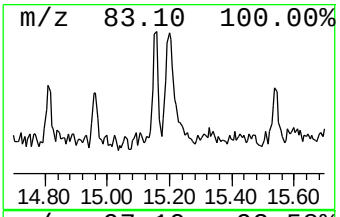
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Carbonic acid, tridecyl 2,2,2-tr...	374	C16H29Cl3O3	1000314-55-9	80
2	Carbonic acid, dodecyl 2,2,2-tri...	360	C15H27Cl3O3	1000314-55-8	80
3	1-Hexadecene, 16-bromo-	302	C16H31Br	118625-56-2	64
4	Carbonic acid, pentadecyl 2,2,2-...	402	C18H33Cl3O3	1000314-56-1	52
5	Disparlure	282	C19H38O	029804-22-6	52

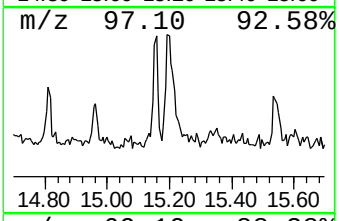


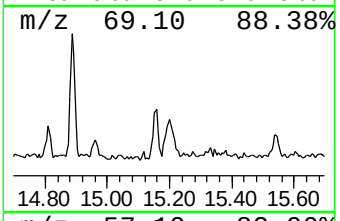


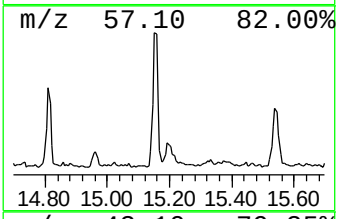


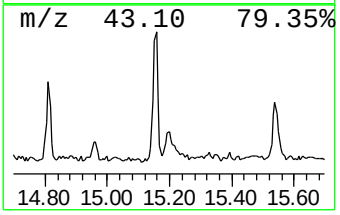












Library Search Compound Report

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Data File : BF118390.D
Acq On : 30 Dec 2019 19:33
Operator : JU
Sample : K6447-01
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : RBR31-48

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF122419.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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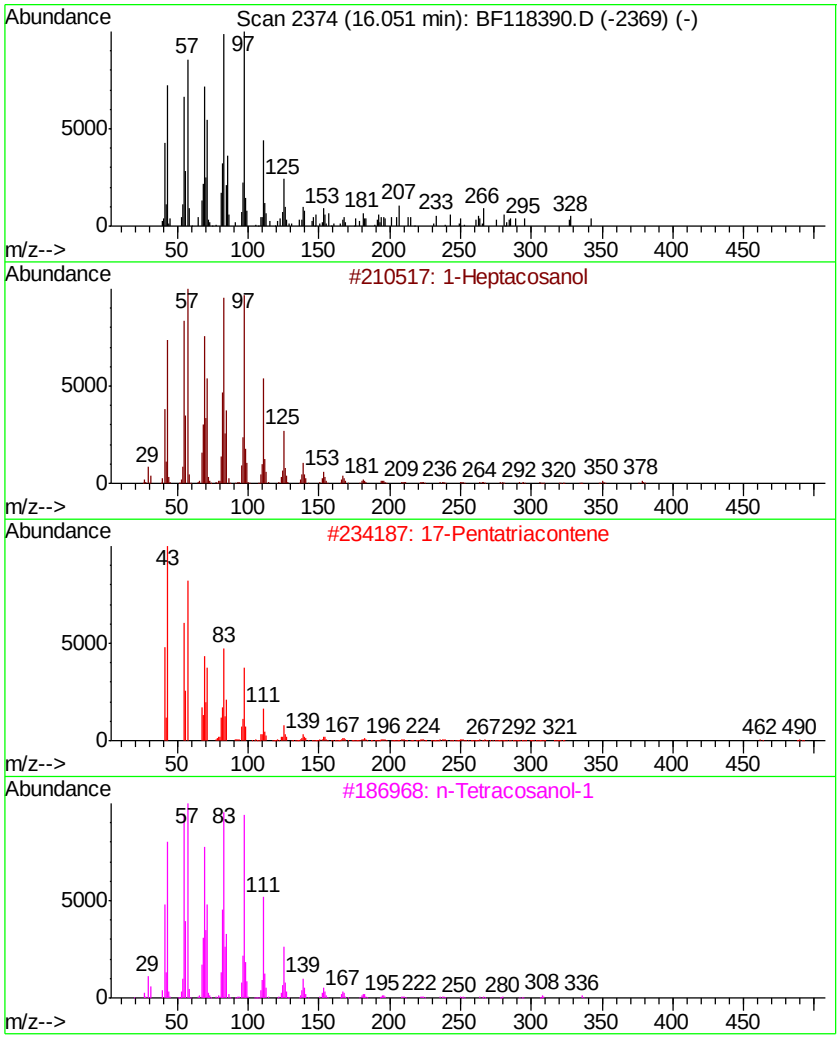
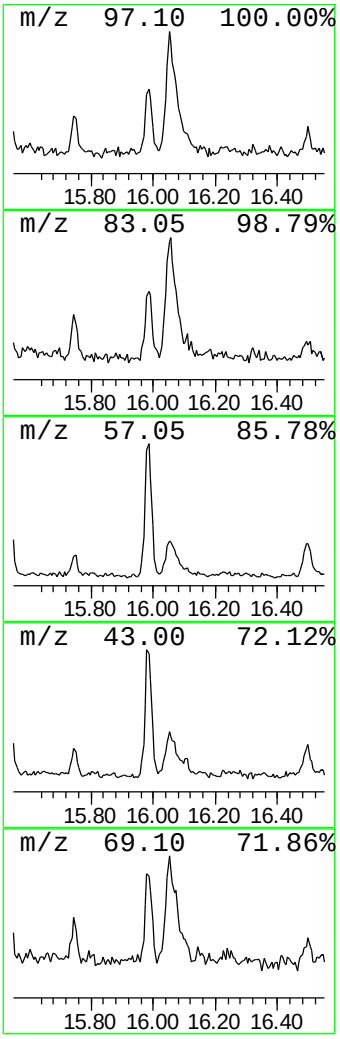
Peak Number 16 1-Heptacosanol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.05	4.85 ng	253742	Perylene-d12	15.54

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heptacosanol	396	C27H56O	002004-39-9	94
2	17-Pentatriacontene	491	C35H70	006971-40-0	91
3	n-Tetracosanol-1	354	C24H50O	000506-51-4	90
4	Methoxyacetic acid, heptadecyl e...	328	C20H40O3	1000282-99-1	87
5	Nonadecyl heptafluorobutyrate	480	C23H39F7O2	1000351-83-9	86





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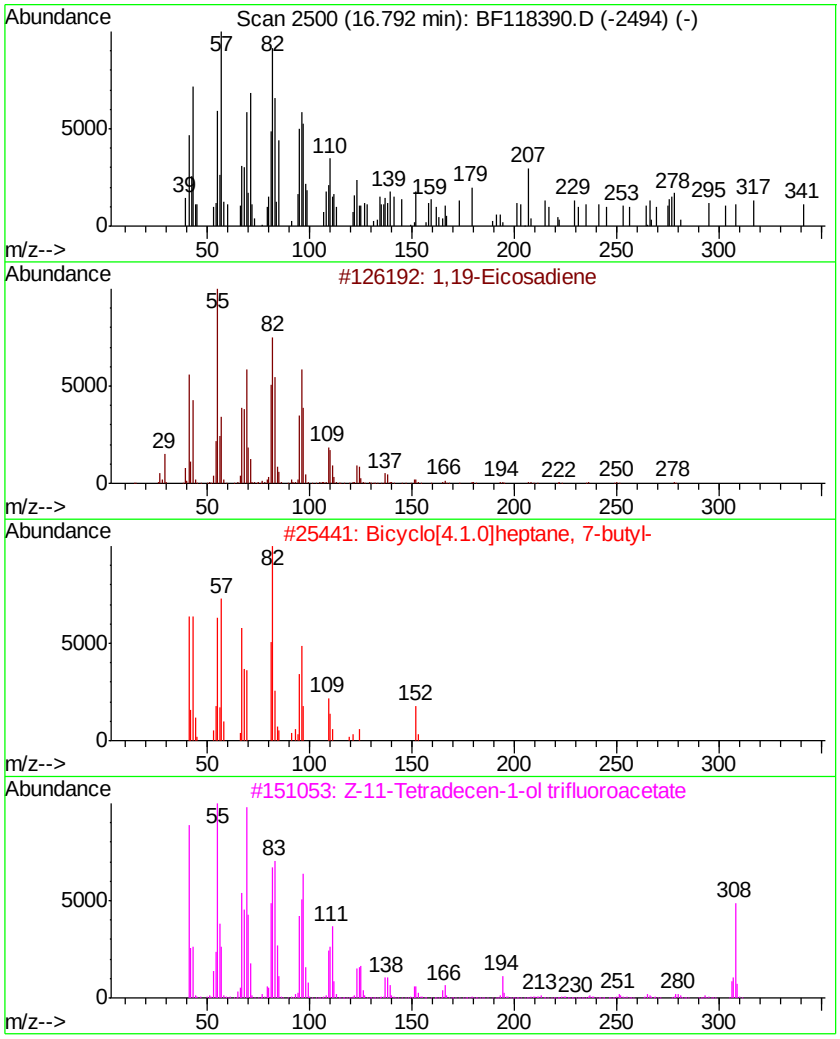
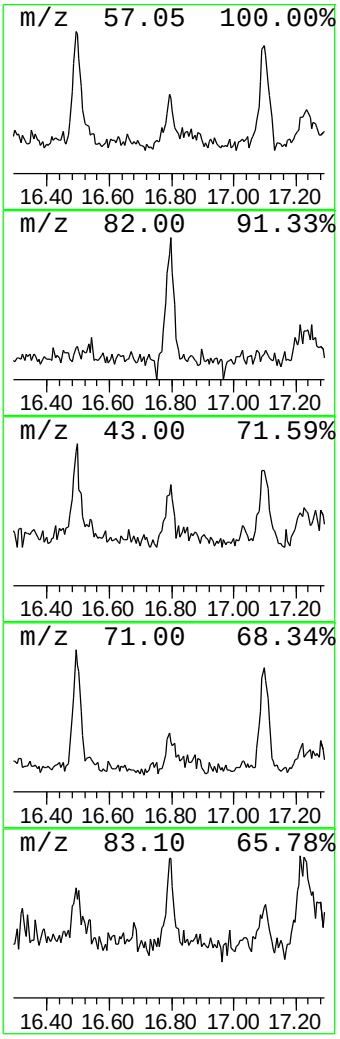
Peak Number 17 1,19-Eicosadiene Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.79	2.04 ng	106638	Perylene-d12	15.54

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,19-Eicosadiene	278	C20H38	014811-95-1	58
2	Bicyclo[4.1.0]heptane, 7-butyl-	152	C11H20	018645-10-8	53
3	Z-11-Tetradecen-1-ol trifluoroac...	308	C16H27F3O2	1000131-33-7	50
4	Bicyclo[10.8.0]eicosane, cis-	278	C20H38	1000155-82-2	49
5	Z,E-3,13-Octadecadien-1-ol	266	C18H34O	1000131-10-4	43





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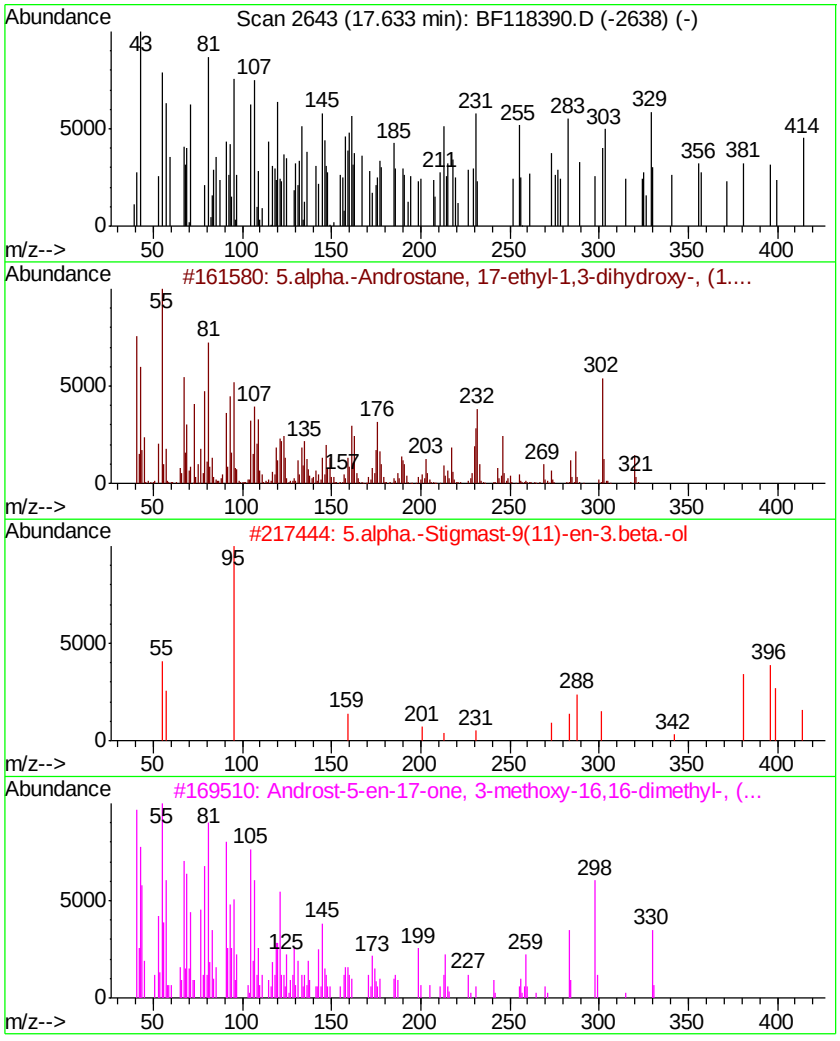
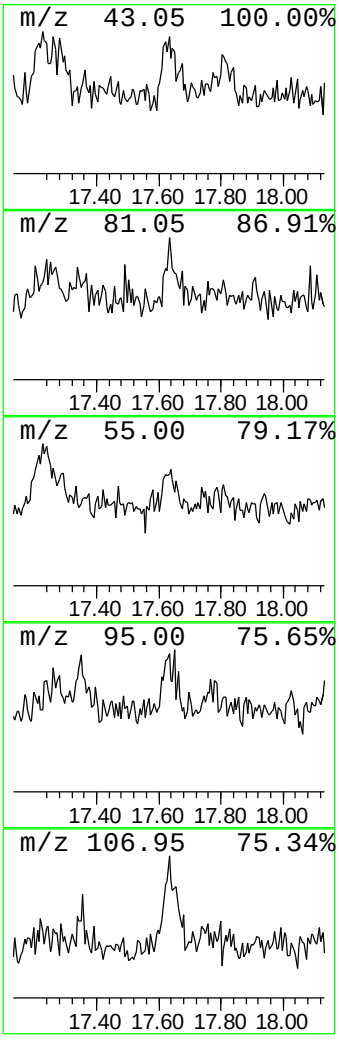
Peak Number 18 unknown17.63 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.63	2.28 ng	119176	Perylene-d12	15.54

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5.alpha.-Androstane, 17-ethyl-1,...	320	C21H36O2	023931-04-6	27
2	5.alpha.-Stigmast-9(11)-en-3.bet...	414	C29H50O	077794-81-1	12
3	Androst-5-en-17-one, 3-methoxy-1...	330	C22H34O2	055837-02-0	12
4	Guanidine, 1,2,3-tri(4-methylphe...	329	C22H23N3	047459-89-2	10
5	Pregn-5-en-3-ol, 21-bromo-20-met...	394	C22H35BrO	055103-80-5	10





Tentatively Identified Compound (LSC) summary

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
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unknown6.62	6.62	79.0	ng	2902310	1	6.85	735133	20.0
Tetradecanoic acid	11.05	3.2	ng	199518	4	11.39	1241760	20.0
n-Hexadecanoic acid	11.93	38.4	ng	2383130	4	11.39	1241760	20.0
Octadecanoic acid	12.71	10.4	ng	646689	4	11.39	1241760	20.0
Cinnamyl cinnamate	13.74	4.5	ng	224477	5	14.05	1000260	20.0
Heptadecyl heptaf...	13.88	14.0	ng	698227	5	14.05	1000260	20.0
Eicosane	14.20	2.1	ng	105498	5	14.05	1000260	20.0
Hexacosane	14.50	3.0	ng	151921	5	14.05	1000260	20.0
Heptadecane	14.81	2.4	ng	123260	6	15.54	1047000	20.0
Heneicosane	15.16	4.5	ng	238106	6	15.54	1047000	20.0
Carbonic acid, tr...	15.20	2.8	ng	147929	6	15.54	1047000	20.0
1-Heptacosanol	16.05	4.8	ng	253742	6	15.54	1047000	20.0
1,19-Eicosadiene	16.79	2.0	ng	106638	6	15.54	1047000	20.0
unknown17.63	17.63	2.3	ng	119176	6	15.54	1047000	20.0