LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120619\
Data File : BP001245.D
Acq On : 06 Dec 2019 22:51
Operator : JU
Sample : K6135-02 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-34

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_P\METHODS\8270-BP112719.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	2.346	41	44	53	rVB2	13499	20918	1.55%	0.111%
2	5.610	592	599	612	rVB	156382	229185	16.96%	1.220%
3	6.210	696	701	715	rVB	769451	988079	73.13%	5.260%
4	7.757	958	964	974	rBV	1004296	1190905	88.14%	6.339%
5	7.945	990	996	1010	rVB	944410	1104116	81.72%	5.877%
6	8.310	1052	1058	1065	rBV	518870	585083	43.30%	3.115%
7	8.551	1094	1099	1104	rVB	763661	880996	65.21%	4.690%
8	9.186	1201	1207	1216	rBV	642150	720768	53.35%	3.837%
9	10.345	1399	1404	1407	rBV	609607	704645	52.15%	3.751%
10	11.116	1528	1535	1540	rBV3	24356	39579	2.93%	0.211%
11	11.392	1578	1582	1585	rBV2	23217	27931	2.07%	0.149%
12	11.504	1598	1601	1606	rVB	27666	34277	2.54%	0.182%
13	11.663	1624	1628	1632	rVB2	22287	27362	2.03%	0.146%
14	11.710	1632	1636	1640	rBV5	15357	20626	1.53%	0.110%
15	12.122	1699	1706	1710	rVB	1152087	1333349	98.69%	7.098%
16	12.251	1723	1728	1729	rBV4	13407	17663	1.31%	0.094%
17	12.328	1737	1741	1745	rBV2	31205	46707	3.46%	0.249%
18	12.422	1754	1757	1765	rVB6	18577	37486	2.77%	0.200%
19	12.545	1773	1778	1782	rVB4	17528	32622	2.41%	0.174%
20	12.657	1791	1797	1801	rBV	24795	45533	3.37%	0.242%
21	12.704	1801	1805	1808	rVV2	21919	30195	2.23%	0.161%
22	12.745	1809	1812	1815	rVV3	18841	27426	2.03%	0.146%
23	12.786	1815	1819	1823	rVV	133164	156495	11.58%	0.833%
24	12.851	1823	1830	1841	rVB4	47350	108696	8.04%	0.579%
25	12.969	1847	1850	1854	rVB	28142	34694	2.57%	0.185%
26	13.204	1884	1890	1895	rBV	737819	892685	66.07%	4.752%
27	13.257	1896	1899	1903	rVV2	34786	48714	3.61%	0.259%
28	13.298	1904	1906	1915	rVB8	18635	30664	2.27%	0.163%
29	13.392	1920	1922	1932	rVB7	11673	27287	2.02%	0.145%
30	13.539	1944	1947	1949	rBV	18824	22792	1.69%	0.121%
31	13.627	1958	1962	1966	rVB4	27955	39806	2.95%	0.212%
32	13.786	1986	1989	1995	rVB4	23047	31732	2.35%	0.169%
33	14.022	2026	2029	2035	rBV	46658	55339	4.10%	0.295%
34	14.104	2040	2043	2046	rVB	27312	30530	2.26%	0.163%

LSC Area Percent Report

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Misc :
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Instrument : BNA_P
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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 >
Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	14.386	2087	2091	2098	rVB3	40172	67889	5.02%	0.361%
36	14.492	2104	2109	2121	rVV	484770	596283	44.13%	3.174%
37	14.598	2124	2127	2134	rVB2	32729	42383	3.14%	0.226%
38	14.804	2157	2162	2164	rBV3	55451	76089	5.63%	0.405%
39	14.833	2164	2167	2170	rVB	53728	67017	4.96%	0.357%
40	14.927	2180	2183	2186	rVB4	19310	21357	1.58%	0.114%
41	15.010	2194	2197	2202	rVB6	16603	19762	1.46%	0.105%
42	15.110	2211	2214	2217	rBV5	18047	28044	2.08%	0.149%
43	15.216	2229	2232	2236	rBV2	18297	26365	1.95%	0.140%
44	15.439	2267	2270	2274	rBV2	18139	20607	1.53%	0.110%
45	15.516	2280	2283	2287	rVB	67061	68568	5.07%	0.365%
46	15.563	2288	2291	2294	rVV	50154	71851	5.32%	0.382%
47	15.604	2294	2298	2301	rVV	682882	821763	60.82%	4.374%
48	15.633	2301	2303	2308	rVB	229734	250228	18.52%	1.332%
49	15.710	2313	2316	2323	rVV	68954	87393	6.47%	0.465%
50	16.157	2389	2392	2395	rBV2	69119	77302	5.72%	0.411%
51	16.298	2412	2416	2420	rBV	227417	234615	17.36%	1.249%
52	16.357	2423	2426	2429	rBV2	47664	58195	4.31%	0.310%
53	16.398	2429	2433	2440	rVB	71972	97820	7.24%	0.521%
54	16.486	2445	2448	2450	rBV2	64156	72597	5.37%	0.386%
55	16.739	2488	2491	2495	rBV	81649	91142	6.75%	0.485%
56	17.139	2556	2559	2564	rBV2	60770	79477	5.88%	0.423%
57	17.257	2575	2579	2581	rBV	485271	556229	41.17%	2.961%
58	17.286	2581	2584	2587	rVV	931768	924526	68.43%	4.921%
59	17.315	2587	2589	2591	rVV2	52441	47077	3.48%	0.251%
60	17.345	2591	2594	2599	rVB	368751	398087	29.46%	2.119%
61	17.410	2602	2605	2608	rBV	103506	110679	8.19%	0.589%
62	17.651	2642	2646	2650	rVB	356757	405334	30.00%	2.158%
63	17.780	2666	2668	2673	rVB	88390	87390	6.47%	0.465%
64	17.880	2681	2685	2689	rVB	1252582	1351116	100.00%	7.192%
65	18.557	2797	2800	2807	rVB	342547	400106	29.61%	2.130%
66	19.121	2893	2896	2903	rVB3	197600	324011	23.98%	1.725%
67	19.245	2912	2917	2920	rBV2	796327	1044755	77.33%	5.561%
68	21.027	3217	3220	3224	rVB	459530	534715	39.58%	2.846%

Sum of corrected areas: 18785657

LSC Report - Integrated Chromatogram

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Misc :
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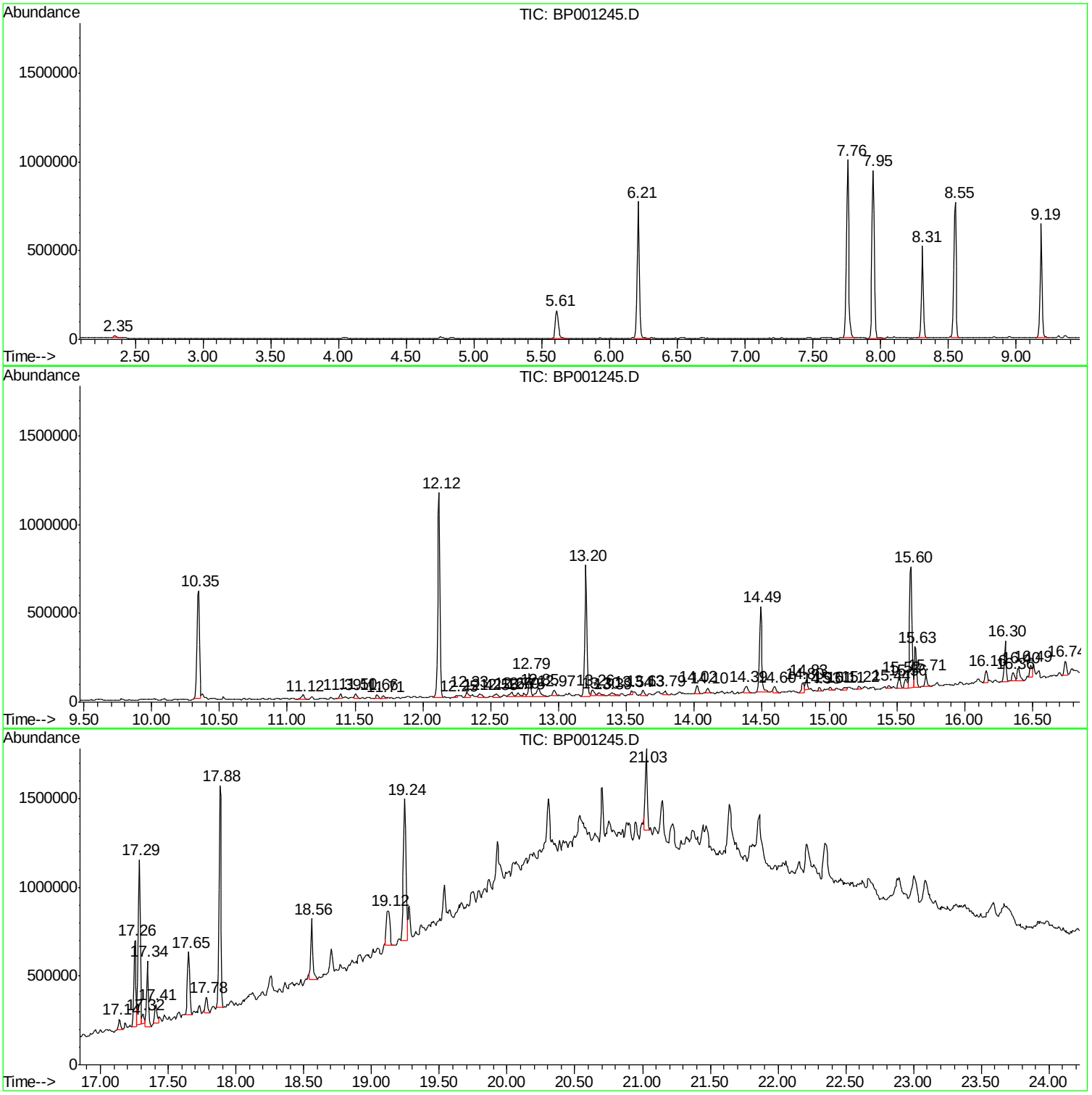
Instrument : BNA_P
ClientSampleId : COMP-34

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.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

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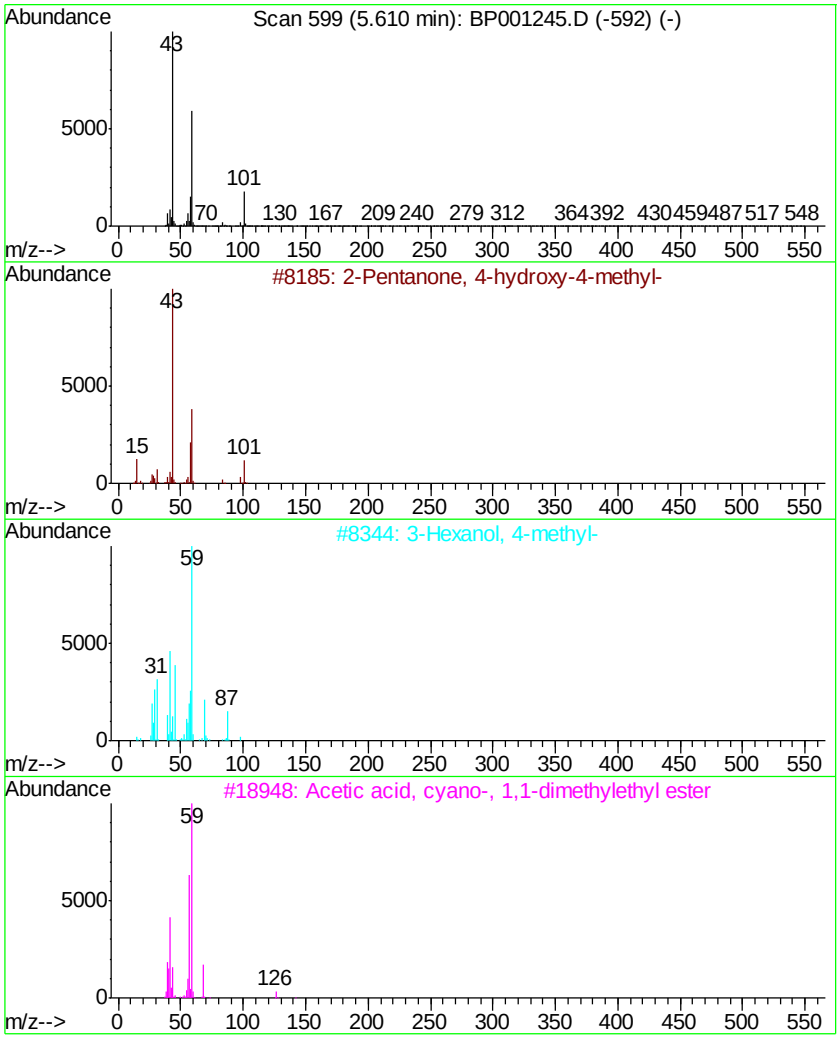
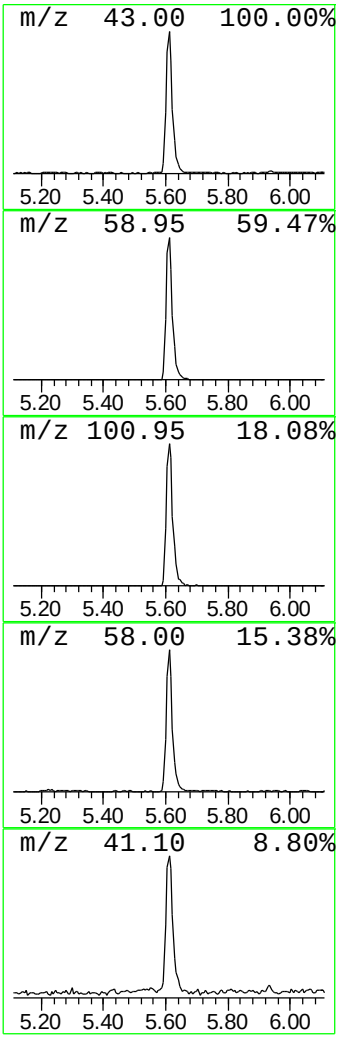
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.61	7.83 ng	229185	1,4-Dichlorobenzene-d4	8.31

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	53
2	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	10
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9





Library Search Compound Report

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Acq On : 06 Dec 2019 22:51
Operator : JU
Sample : K6135-02 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-34

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

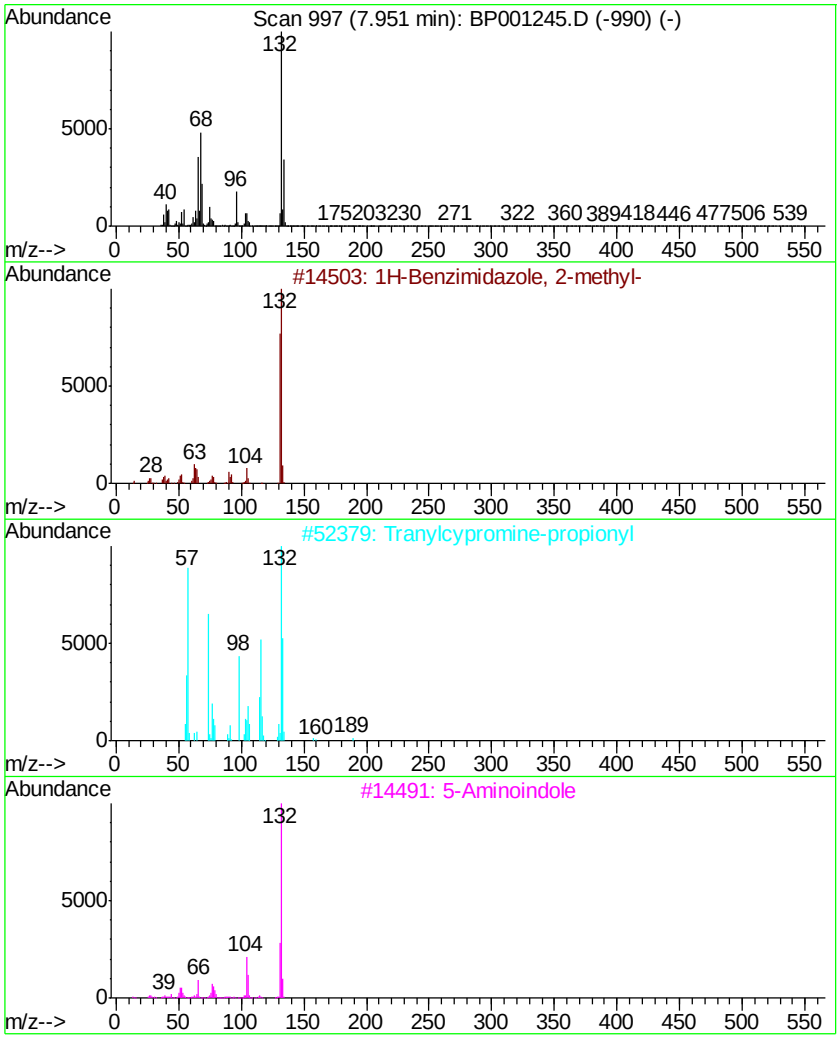
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

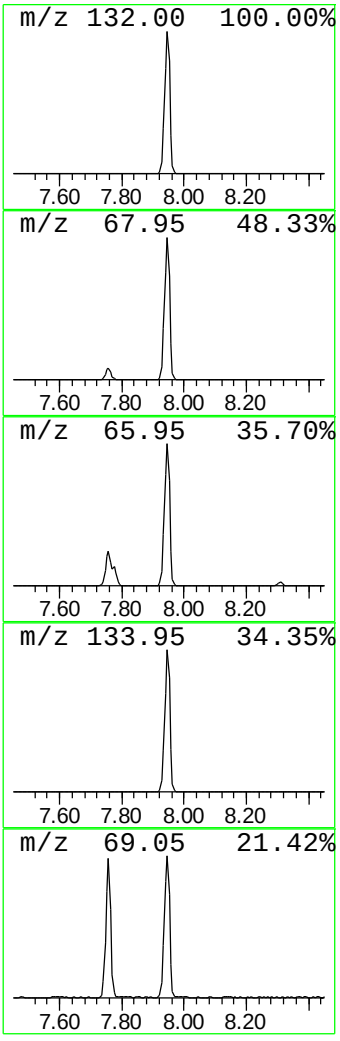
****

Peak Number 2 unknown7.95 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.95	37.74 ng	1104120	1,4-Dichlorobenzene-d4	8.31

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22
2	Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	22
3	5-Aminoindole	132	C8H8N2	005192-03-0	22
4	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
5	Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	14





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Sample : K6135-02 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-34

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

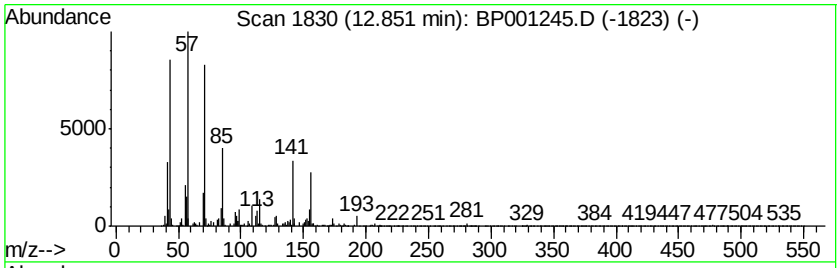
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

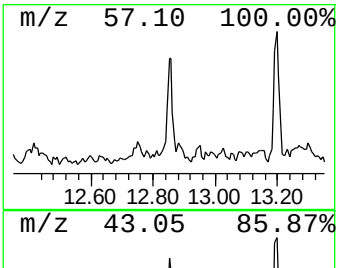
****

Peak Number 4 Dodecane, 4,6-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.85	2.44 ng	108696	Acenaphthene-d10	13.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	55
2	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	53
3	1,3-Dimethylcyclopentanol	114	C7H14O	019550-46-0	49
4	Heptadecane, 2,6,10,14-tetramethyl-	296	C21H44	018344-37-1	49
5	Undecane, 6-ethyl-	184	C13H28	017312-60-6	47





Library Search Compound Report

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Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-34

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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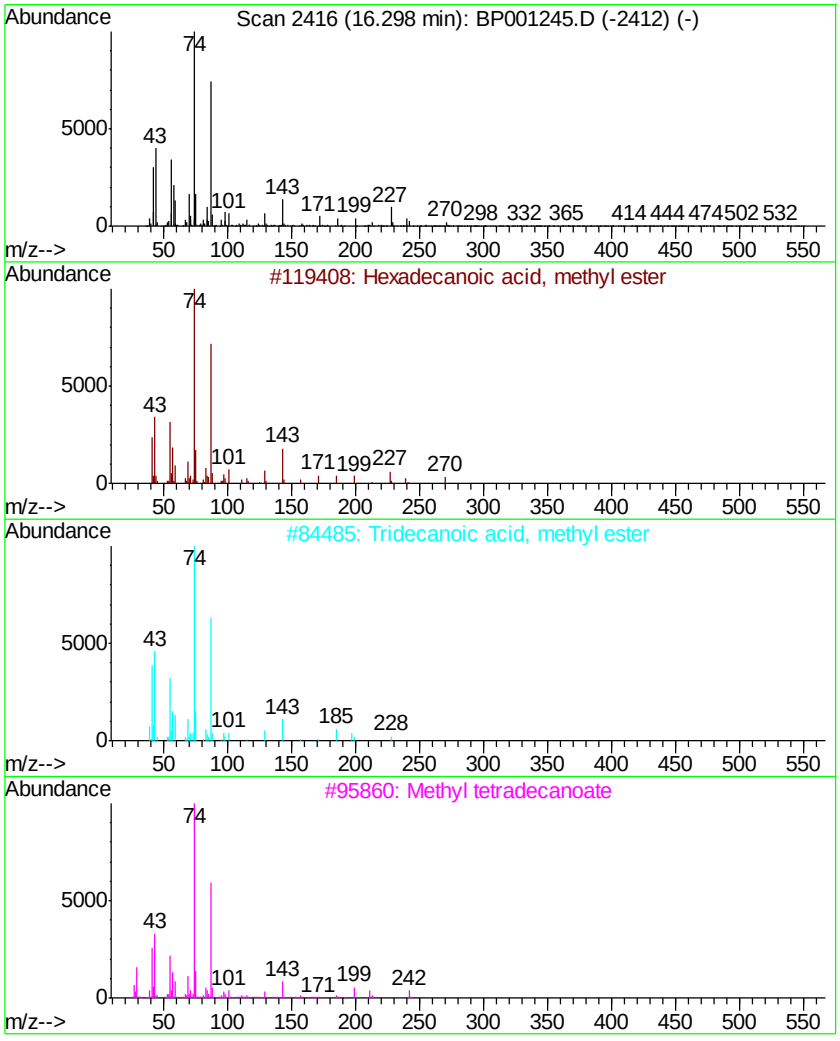
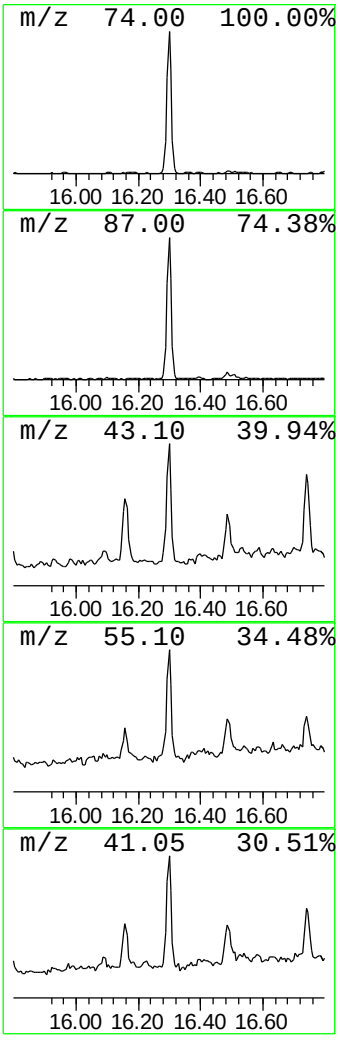
****

Peak Number 5 Hexadecanoic acid, methyl e... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.30	5.71 ng	234615	Phenanthrene-d10	15.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	95
2	Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	91
3	Methyl tetradecanoate	242	C15H30O2	000124-10-7	87
4	Hexadecanoic acid, 15-methyl-, m...	284	C18H36O2	006929-04-0	86
5	Dodecanoic acid, 10-methyl-, met...	228	C14H28O2	005129-65-7	86





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Instrument : BNA_P
ClientSampleId : COMP-34

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

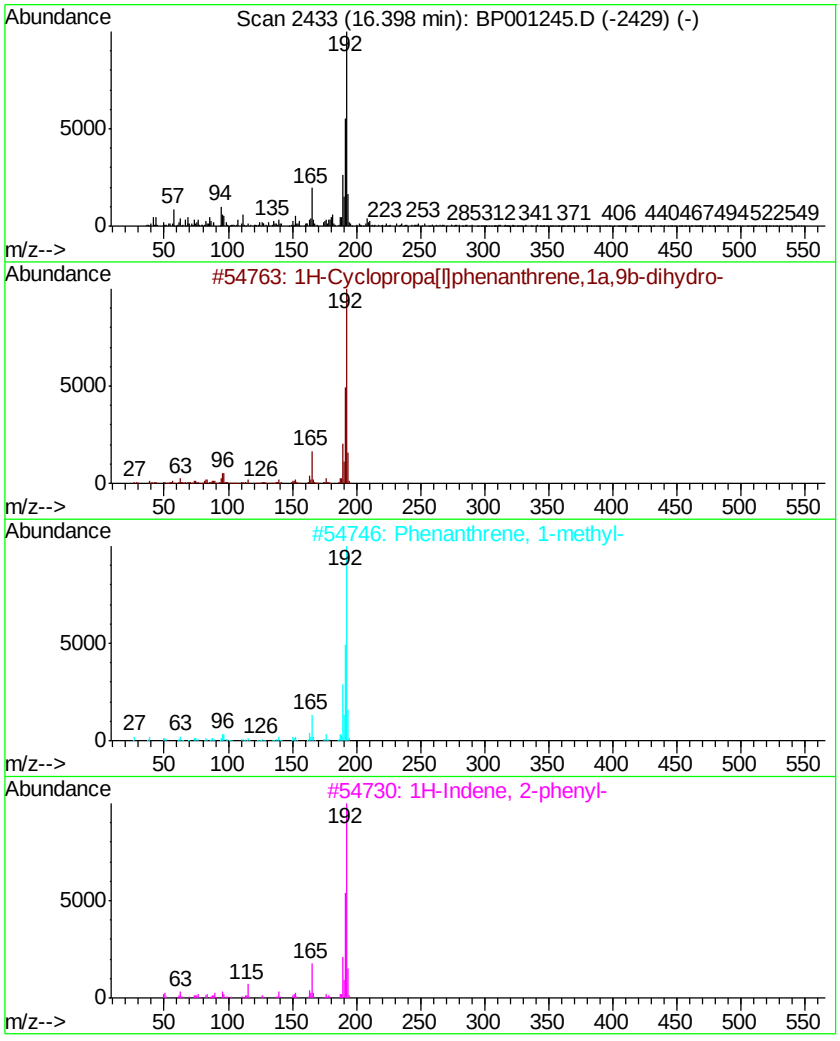
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

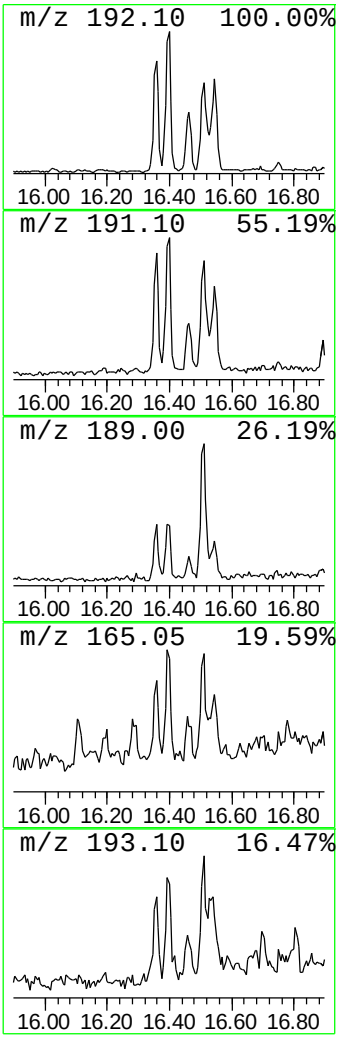
****

Peak Number 6 1H-Cyclopropa[l]phenanthren... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.40	2.38 ng	97820	Phenanthrene-d10	15.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	96
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
3	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	94
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	94





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Misc :
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Instrument : BNA_P
ClientSampleId : COMP-34

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

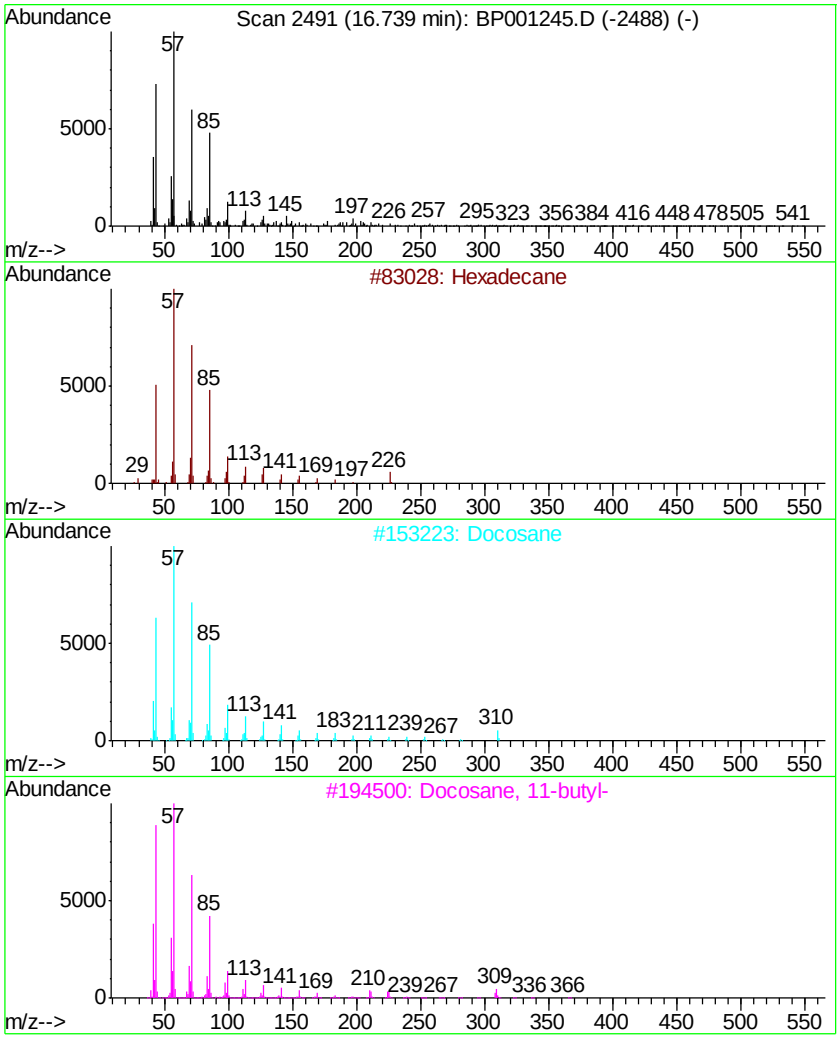
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

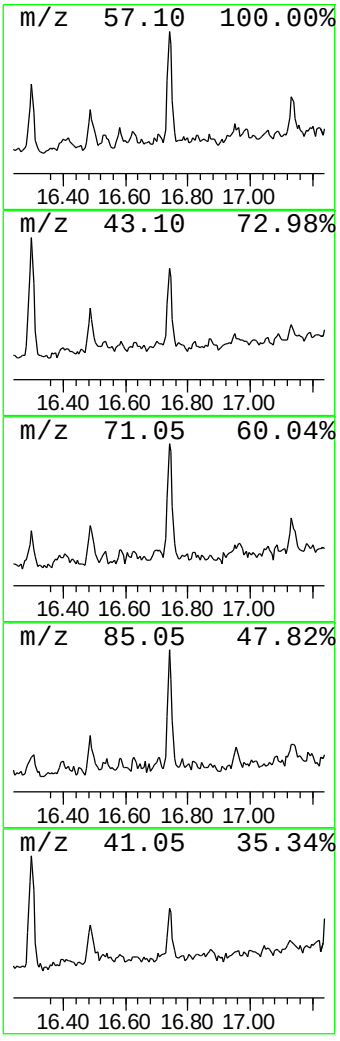
****

Peak Number 7 Hexadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.74	2.22 ng	91142	Phenanthrene-d10	15.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	90
2	Docosane	310	C22H46	000629-97-0	87
3	Docosane, 11-butyl-	366	C26H54	013475-76-8	83
4	Heneicosane	296	C21H44	000629-94-7	83
5	Hexacosane	366	C26H54	000630-01-3	83





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Data File : BP001245.D
Acq On : 06 Dec 2019 22:51
Operator : JU
Sample : K6135-02 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-34

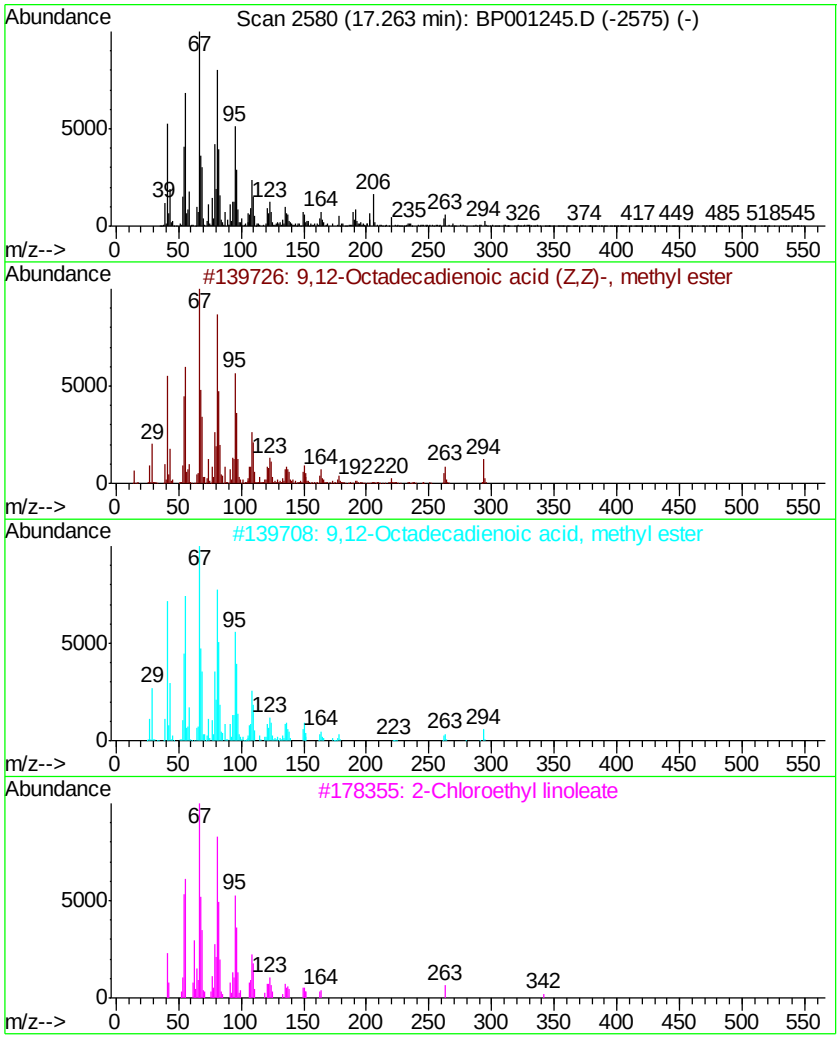
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

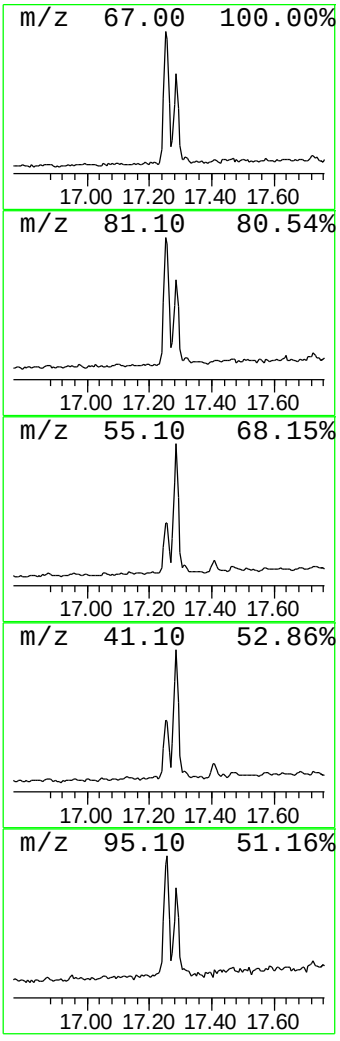
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 9,12-Octadecadienoic acid (... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.26	13.54 ng	556229	Phenanthrene-d10	15.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
2	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99
3	2-Chloroethyl linoleate	342	C20H35ClO2	025525-76-2	99
4	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
5	10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	98





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120619\
Data File : BP001245.D
Acq On : 06 Dec 2019 22:51
Operator : JU
Sample : K6135-02 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-34

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.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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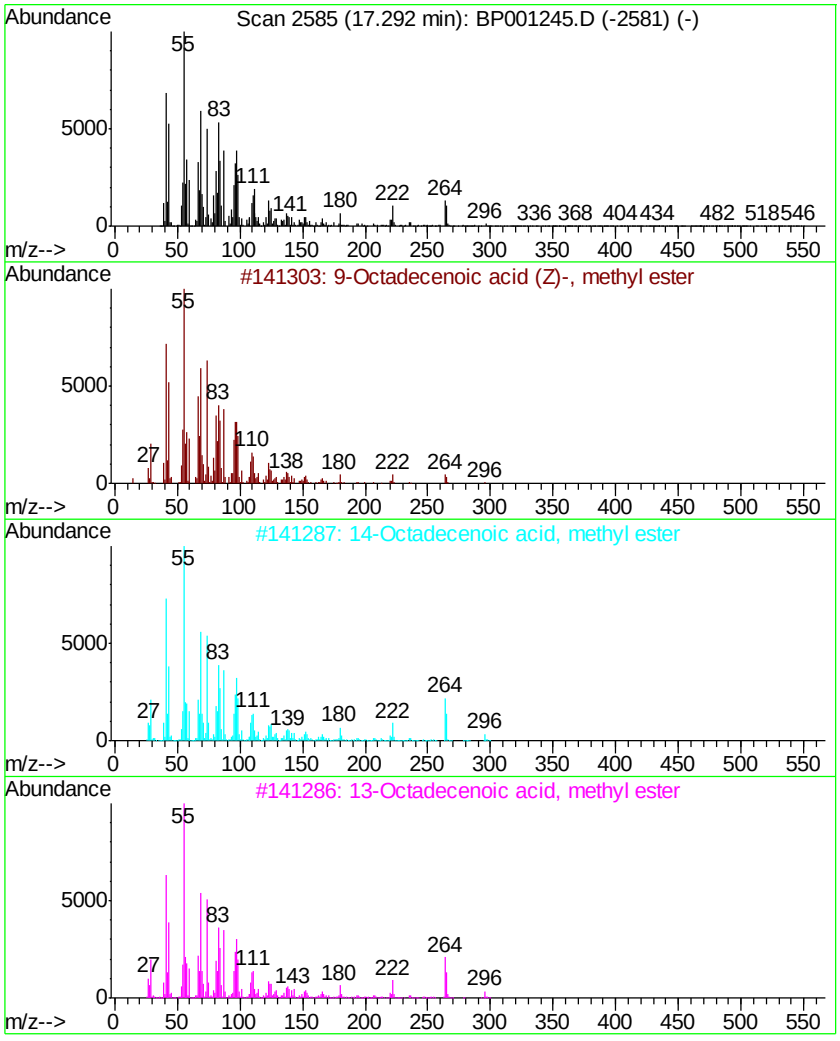
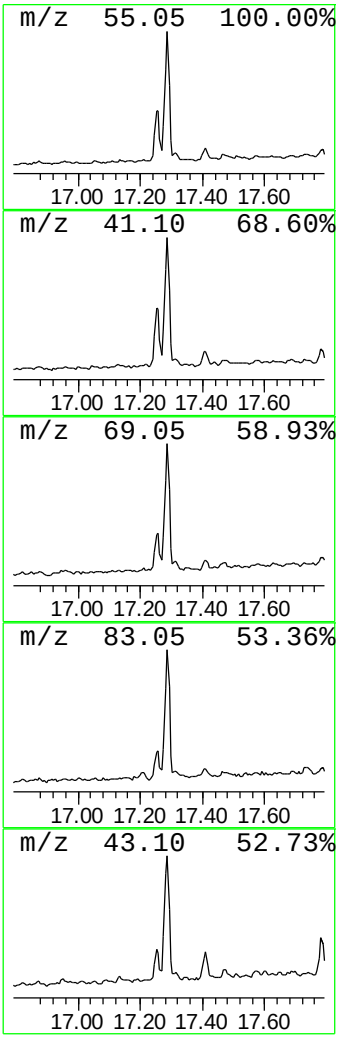
Peak Number 9 9-Octadecenoic acid (Z)-, m... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.29	22.50 ng	924526	Phenanthrene-d10	15.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	98
2	14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	98
3	13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	98
4	9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	96
5	10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	94





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Instrument : BNA_P
ClientSampleId : COMP-34

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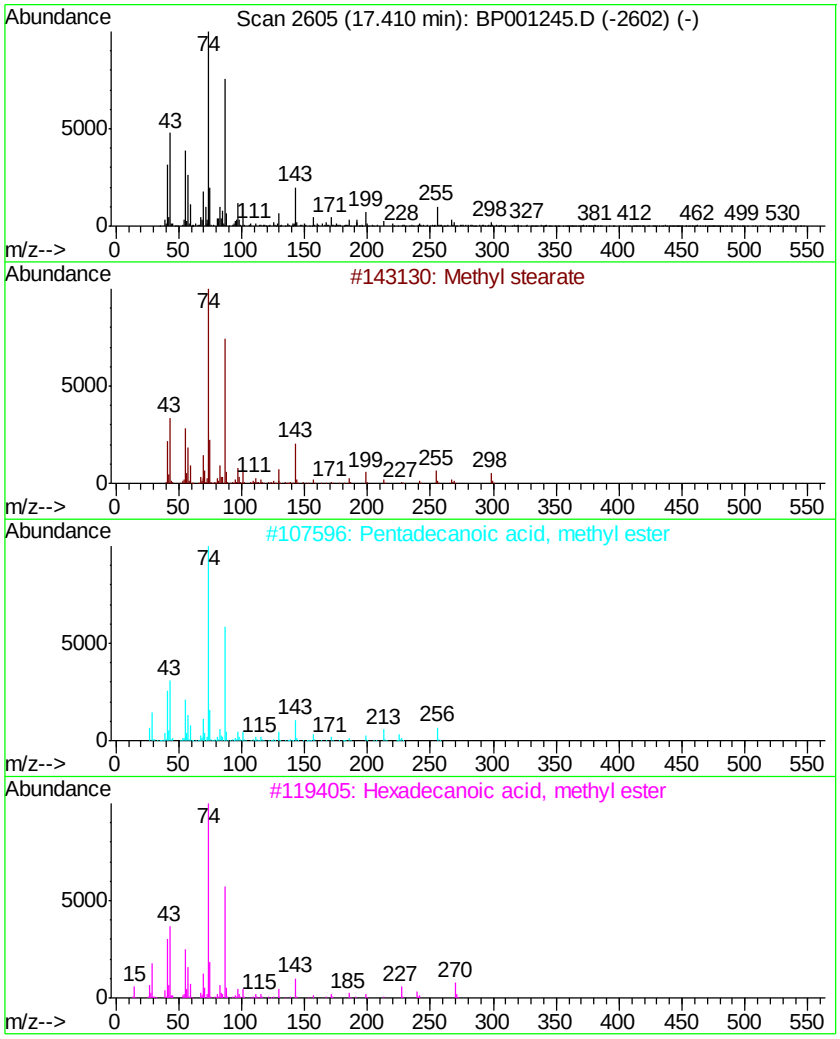
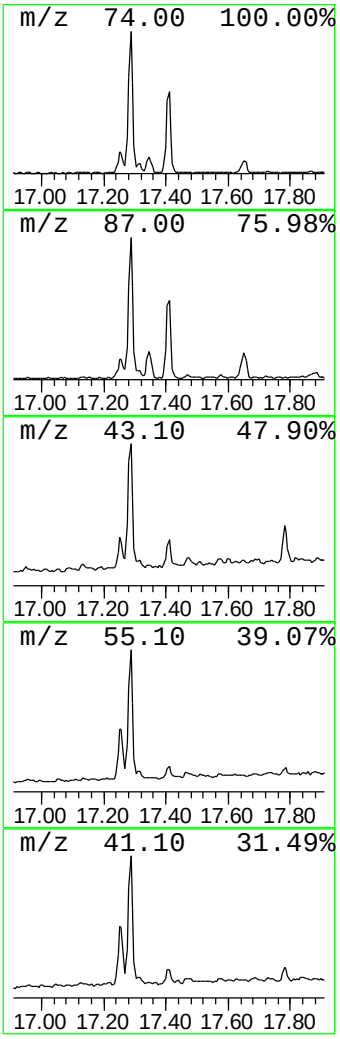
Peak Number 10 Pentadecanoic acid, methyl ... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.41	2.69 ng	110679	Phenanthrene-d10	15.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl stearate	298	C19H38O2	000112-61-8	91
2	Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	87
3	Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	80
4	Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	80
5	Methyl 11-methyl-dodecanoate	228	C14H28O2	1000336-45-1	80





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Operator : JU
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Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-34

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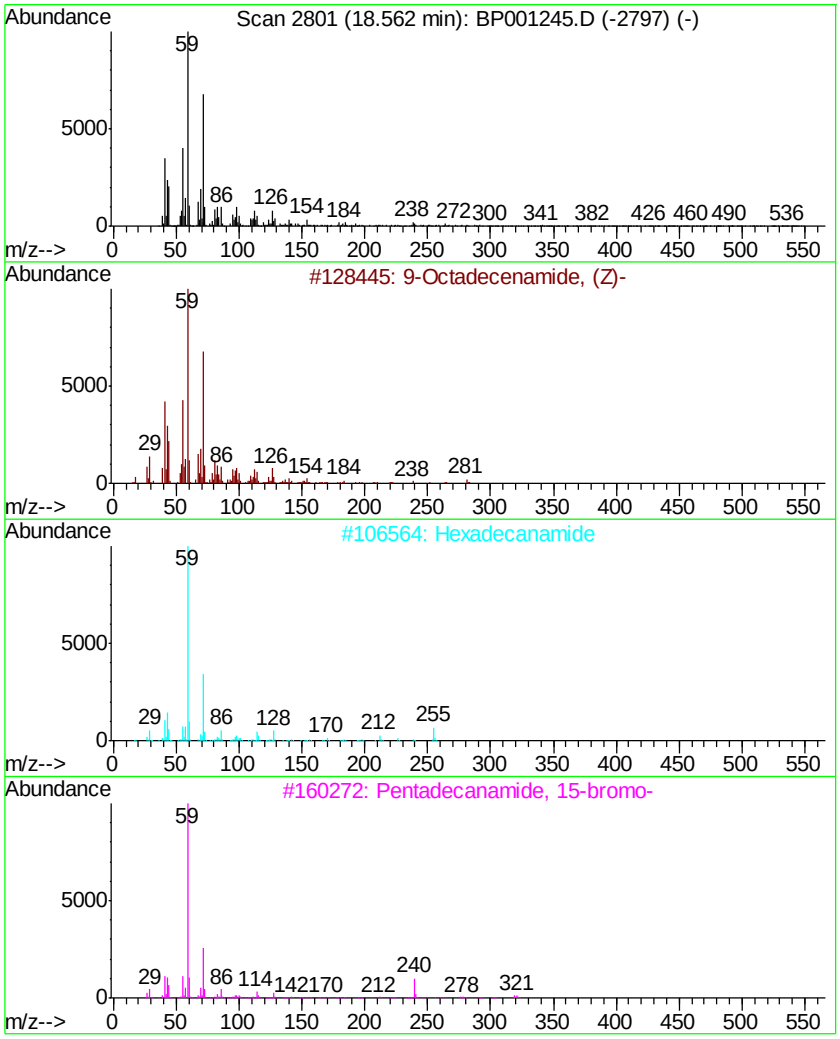
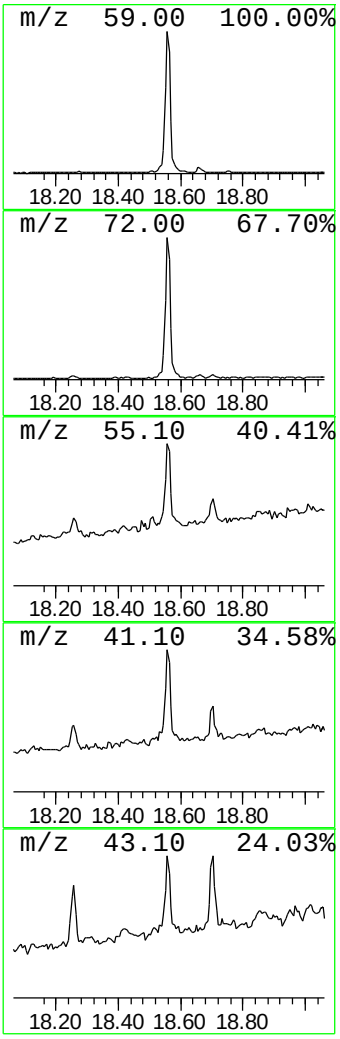
Peak Number 11 9-Octadecenamide, (Z)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.56	7.66 ng	400106	Chrysene-d12	19.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	94
2	Hexadecanamide	255	C16H33NO	000629-54-9	59
3	Pentadecanamide, 15-bromo-	319	C15H30BrNO	1000163-86-1	59
4	Pentanamide, 4-methyl-	115	C6H13NO	001119-29-5	50
5	Tetradecanamide	227	C14H29NO	000638-58-4	45





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Sample : K6135-02 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-34

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TIC Integration Parameters: LSCINT.P

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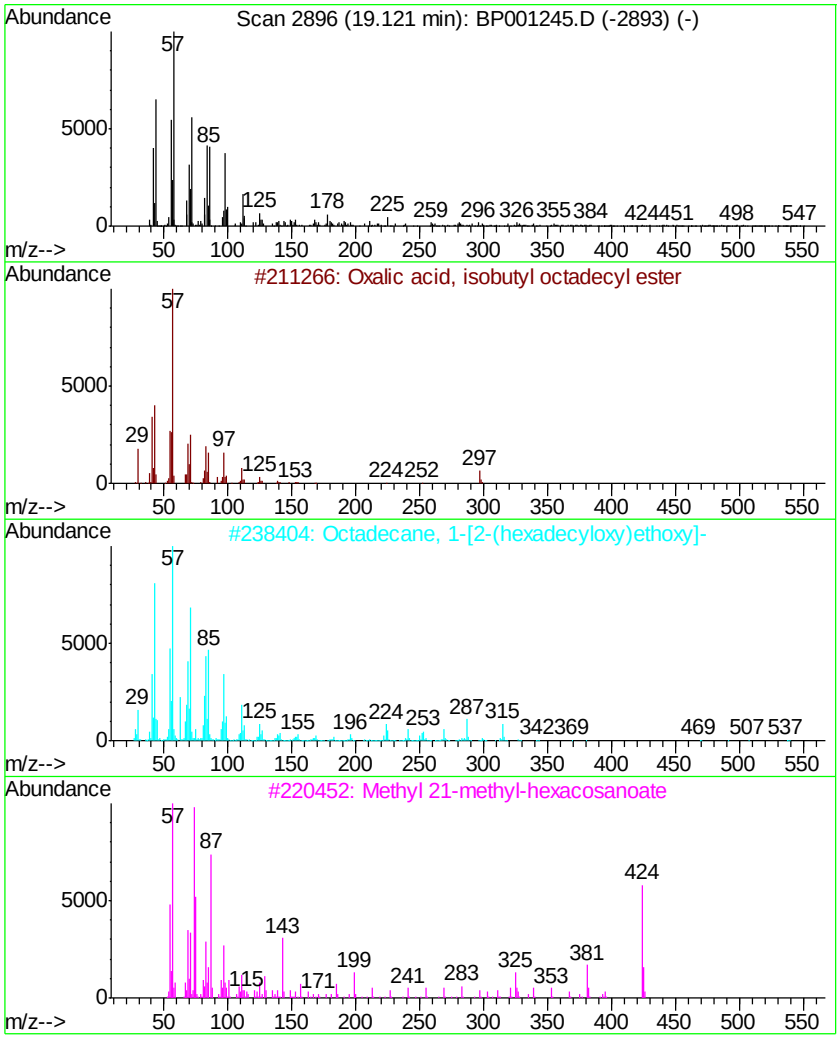
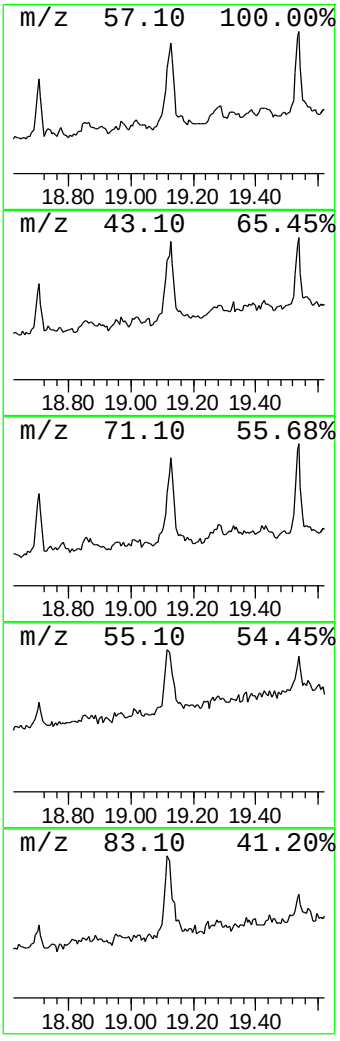
Peak Number 12 Oxalic acid, isobutyl octad... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.12	6.20 ng	324011	Chrysene-d12	19.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxalic acid, isobutyl octadecyl ...	398	C24H46O4	1000309-38-3	91
2	Octadecane, 1-[2-(hexadecyloxy)e...	539	C36H74O2	017367-10-1	87
3	Methyl 21-methyl-hexacosanoate	424	C28H56O2	1000336-27-5	83
4	Cyclotetradecane	196	C14H28	000295-17-0	83
5	Oxalic acid, isobutyl hexadecyl ...	370	C22H42O4	1000309-38-1	83





Tentatively Identified Compound (LSC) summary

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Data File : BP001245.D
Acq On : 06 Dec 2019 22:51
Operator : JU
Sample : K6135-02 2X
Misc :
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ClientSampleId : COMP-34

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.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
2-Pentanone, 4-hy...	5.61	7.8	ng	229185	1	8.31	585083	20.0
unknown7.95	7.95	37.7	ng	1104120	1	8.31	585083	20.0
Dodecane, 4,6-dim...	12.85	2.4	ng	108696	3	13.20	892685	20.0
Hexadecanoic acid...	16.30	5.7	ng	234615	4	15.60	821763	20.0
1H-Cyclopropa[l]p...	16.40	2.4	ng	97820	4	15.60	821763	20.0
Hexadecane	16.74	2.2	ng	91142	4	15.60	821763	20.0
9,12-Octadecadien...	17.26	13.5	ng	556229	4	15.60	821763	20.0
9-Octadecenoic ac...	17.29	22.5	ng	924526	4	15.60	821763	20.0
Pentadecanoic aci...	17.41	2.7	ng	110679	4	15.60	821763	20.0
9-Octadecenamide,...	18.56	7.7	ng	400106	5	19.24	1044760	20.0
Oxalic acid, isob...	19.12	6.2	ng	324011	5	19.24	1044760	20.0