LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120619\
Data File : BP001243.D
Acq On : 06 Dec 2019 21:51
Operator : JU
Sample : K6135-06 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-36

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	55	rVB	32628	48997	3.42%	0.255%
2	5.610	595	599	609	rVB	154198	220078	15.35%	1.147%
3	6.210	695	701	713	rVB	862181	1092660	76.22%	5.694%
4	7.757	958	964	978	rBV	1013885	1254205	87.49%	6.536%
5	7.946	990	996	1009	rVB	1016663	1181283	82.40%	6.156%
6	8.310	1052	1058	1066	rVB	483383	562071	39.21%	2.929%
7	8.545	1093	1098	1104	rBV	770408	904087	63.07%	4.711%
8	9.187	1201	1207	1215	rBV	650717	741221	51.71%	3.862%
9	9.945	1331	1336	1342	rBV	93469	109256	7.62%	0.569%
10	10.340	1398	1403	1407	rBV	598687	696207	48.57%	3.628%
11	10.369	1407	1408	1415	rVB2	17455	19747	1.38%	0.103%
12	10.528	1431	1435	1440	rBV	17968	20306	1.42%	0.106%
13	10.975	1507	1511	1517	rBV8	7878	15242	1.06%	0.079%
14	11.116	1529	1535	1541	rBV3	26559	41004	2.86%	0.214%
15	11.392	1574	1582	1586	rBV6	19029	36102	2.52%	0.188%
16	11.504	1597	1601	1605	rVB3	15604	19621	1.37%	0.102%
17	11.663	1624	1628	1633	rVB7	13306	17040	1.19%	0.089%
18	11.710	1633	1636	1640	rBV3	12786	17544	1.22%	0.091%
19	12.122	1699	1706	1710	rBV	1156965	1413656	98.61%	7.366%
20	12.328	1737	1741	1745	rBV4	16868	31427	2.19%	0.164%
21	12.745	1809	1812	1815	rBV2	18355	21498	1.50%	0.112%
22	12.786	1815	1819	1823	rVV	118484	120131	8.38%	0.626%
23	12.851	1827	1830	1834	rVB2	39426	54284	3.79%	0.283%
24	12.969	1846	1850	1853	rBV	25545	29173	2.04%	0.152%
25	13.081	1867	1869	1873	rVB5	19254	18696	1.30%	0.097%
26	13.204	1884	1890	1895	rBV	715433	860446	60.02%	4.484%
27	13.257	1896	1899	1902	rVB	26835	33010	2.30%	0.172%
28	13.539	1944	1947	1950	rBV	20416	30295	2.11%	0.158%
29	13.622	1958	1961	1967	rVB3	27760	40164	2.80%	0.209%
30	13.733	1977	1980	1985	rBV4	14312	27321	1.91%	0.142%
31	13.786	1986	1989	1995	rVB4	20518	27162	1.89%	0.142%
32	14.022	2026	2029	2034	rBV3	25111	34264	2.39%	0.179%
33	14.104	2039	2043	2051	rVB	41176	63904	4.46%	0.333%
34	14.386	2086	2091	2099	rVB3	50861	96491	6.73%	0.503%

LSC Area Percent Report

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Sample : K6135-06 2X
Misc :
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Instrument : BNA_P
ClientSampleId : COMP-36

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	14.492	2104	2109	2122	rVB	761460	929512	64.84%	4.844%
36	14.828	2160	2166	2170	rBV	80002	121882	8.50%	0.635%
37	14.922	2179	2182	2186	rBV3	20919	28662	2.00%	0.149%
38	15.010	2194	2197	2201	rVB5	18446	21933	1.53%	0.114%
39	15.104	2210	2213	2216	rBV4	16377	27445	1.91%	0.143%
40	15.216	2228	2232	2237	rBV	30061	53286	3.72%	0.278%
41	15.433	2267	2269	2273	rBV	23641	24750	1.73%	0.129%
42	15.516	2280	2283	2287	rBV	26119	29829	2.08%	0.155%
43	15.563	2287	2291	2293	rBV	62577	76868	5.36%	0.401%
44	15.598	2293	2297	2300	rVV	692973	820437	57.23%	4.275%
45	15.633	2300	2303	2309	rVB	482697	525233	36.64%	2.737%
46	15.686	2309	2312	2313	rBV3	14036	15709	1.10%	0.082%
47	15.710	2313	2316	2322	rVB	126005	140846	9.82%	0.734%
48	15.957	2355	2358	2363	rVB	48238	55689	3.88%	0.290%
49	16.157	2389	2392	2395	rBV	27482	26884	1.88%	0.140%
50	16.357	2422	2426	2429	rBV	55201	66993	4.67%	0.349%
51	16.392	2429	2432	2440	rVB	89938	110567	7.71%	0.576%
52	16.463	2440	2444	2445	rBV	38119	41316	2.88%	0.215%
53	16.486	2445	2448	2449	rBV	50464	49887	3.48%	0.260%
54	16.504	2449	2451	2455	rVB2	83661	85298	5.95%	0.444%
55	16.739	2489	2491	2496	rVB2	27687	28497	1.99%	0.148%
56	16.786	2496	2499	2509	rBV2	57222	102157	7.13%	0.532%
57	17.139	2556	2559	2563	rVB2	44085	52661	3.67%	0.274%
58	17.180	2564	2566	2570	rBV4	24461	31138	2.17%	0.162%
59	17.345	2590	2594	2598	rBV	686702	715006	49.88%	3.726%
60	17.651	2641	2646	2650	rVB	615381	712773	49.72%	3.714%
61	17.880	2680	2685	2689	rVB	1400158	1433552	100.00%	7.470%
62	18.557	2797	2800	2807	rVB	198387	206865	14.43%	1.078%
63	19.116	2892	2895	2902	rBV2	163206	273885	19.11%	1.427%
64	19.245	2911	2917	2920	rBV2	731171	1142357	79.69%	5.953%
65	19.274	2920	2922	2926	rVB	260416	244201	17.03%	1.273%
66	20.527	3132	3135	3139	rBV	187893	281576	19.64%	1.467%
67	20.939	3203	3205	3209	rVB	181045	200256	13.97%	1.044%
68	21.015	3214	3218	3222	rVV2	470080	614095	42.84%	3.200%

Sum of corrected areas: 19190638

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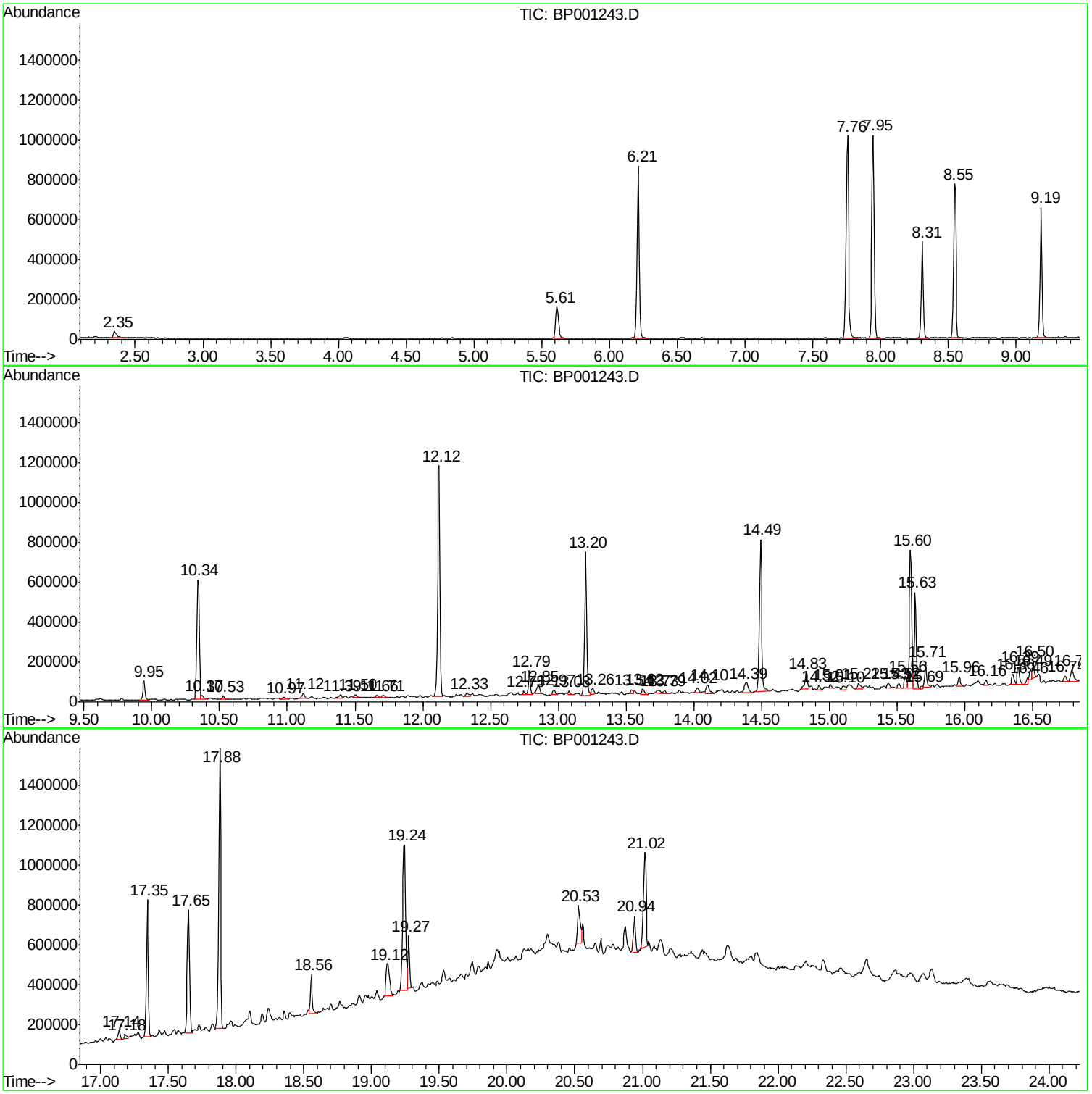
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
COMP-36

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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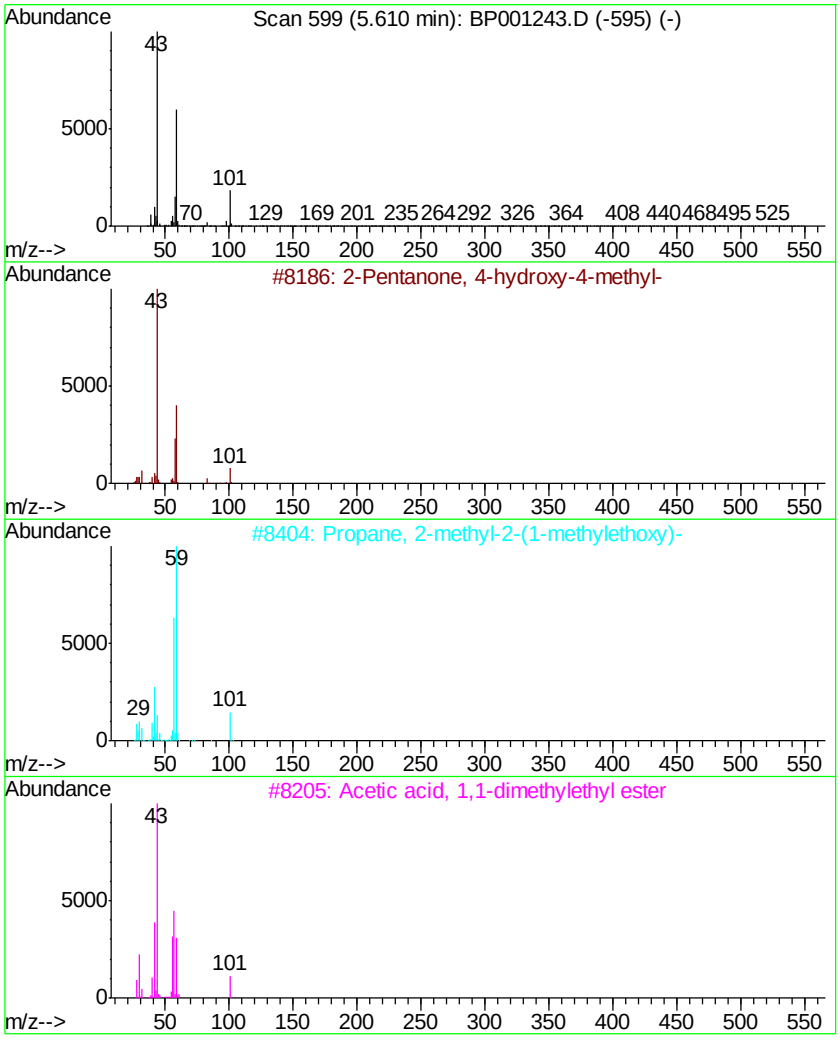
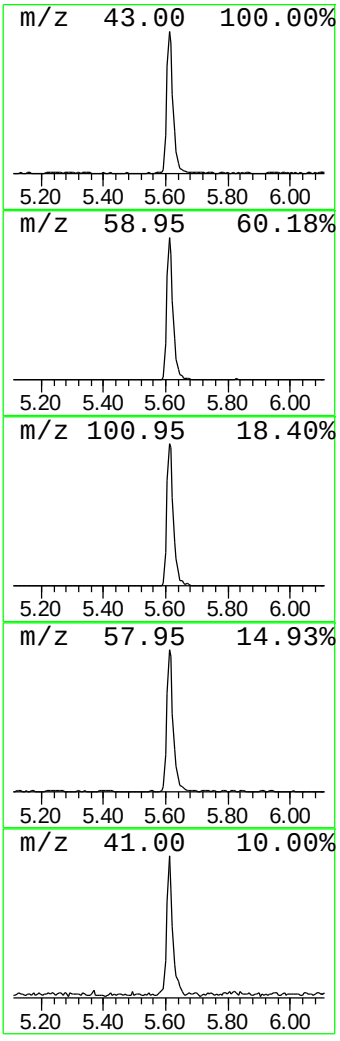
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.61	7.83 ng	220078	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	50
2	Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
3	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	28
4	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
5	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120619\
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Operator : JU
Sample : K6135-06 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-36

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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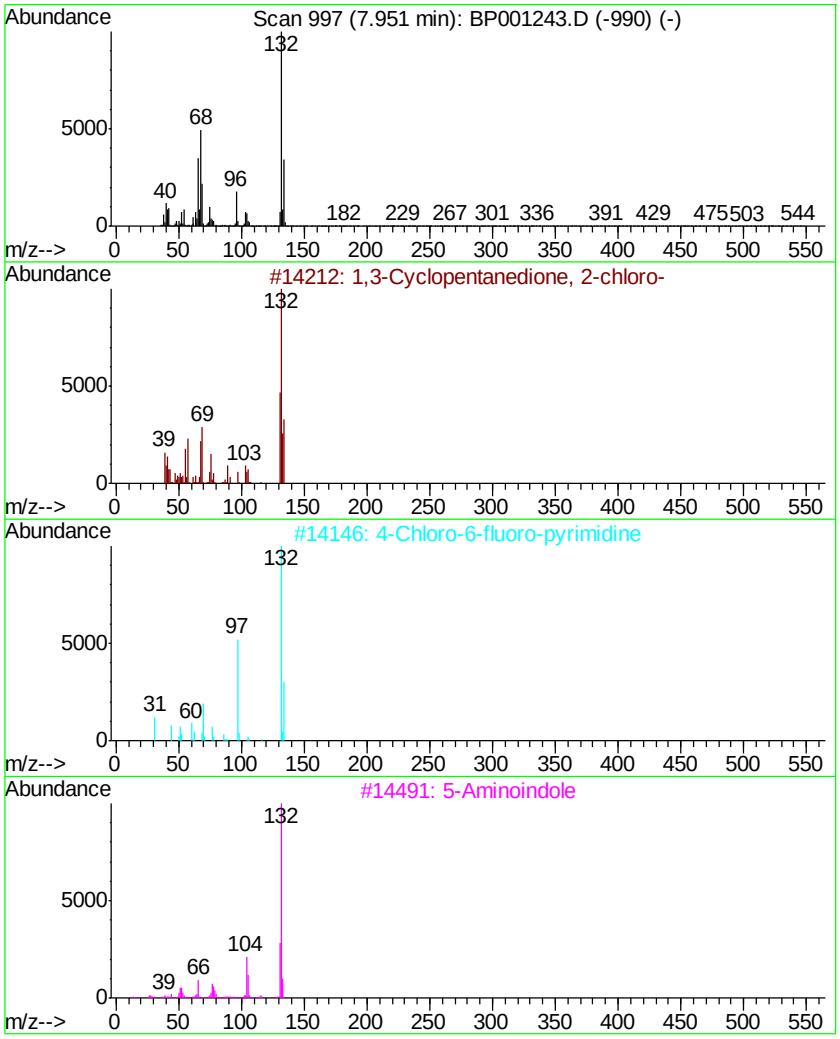
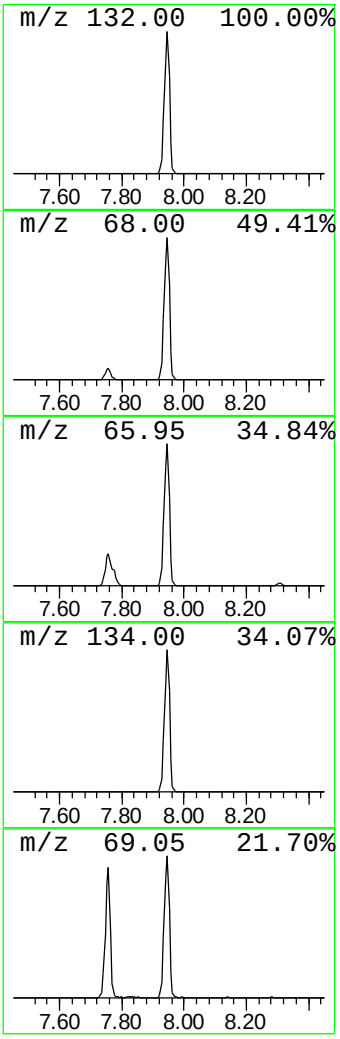
****

Peak Number 2 unknown7.95 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.95	42.03 ng	1181280	1,4-Dichlorobenzene-d4	8.31

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	28
2	4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	27
3	5-Aminoindole	132	C8H8N2	005192-03-0	22
4	Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	22
5	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	22





Library Search Compound Report

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Sample : K6135-06 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-36

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.M
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TIC Library : C:\DATABASE\NIST11.L
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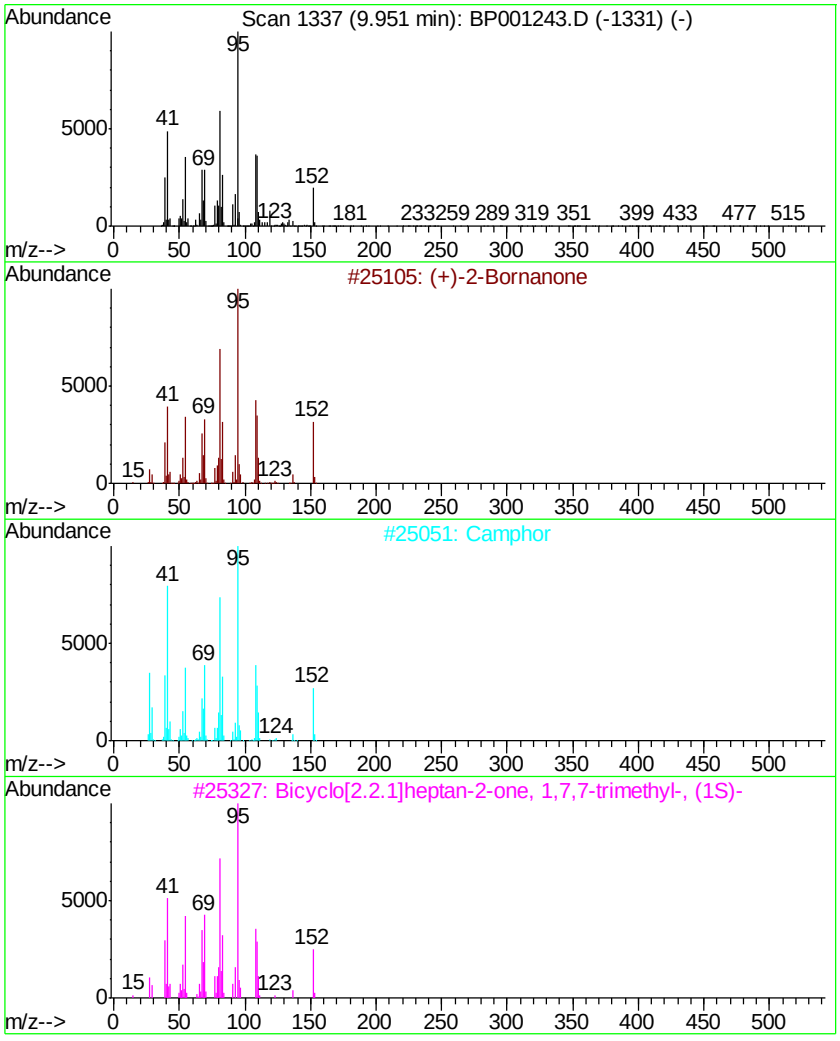
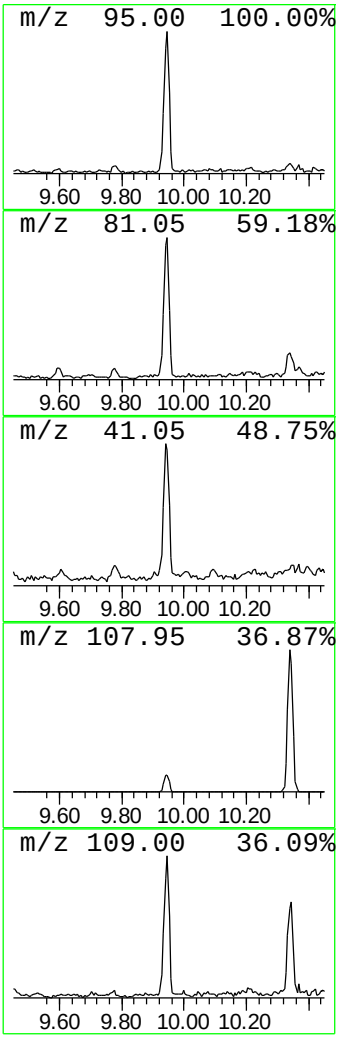
****

Peak Number 4 (+)-2-Bornanone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.95	3.14 ng	109256	Naphthalene-d8	10.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	(+)-2-Bornanone	152	C10H16O	000464-49-3	96
2	Camphor	152	C10H16O	000076-22-2	94
3	Bicyclo[2.2.1]heptan-2-one, 1,7,...	152	C10H16O	000464-48-2	93
4	5,7-Octadien-4-one, 2,6-dimethyl...	152	C10H16O	006752-80-3	52
5	Bicyclo[3.1.0]hexane, 1,5-dimethyl-	110	C8H14	1000142-17-5	46





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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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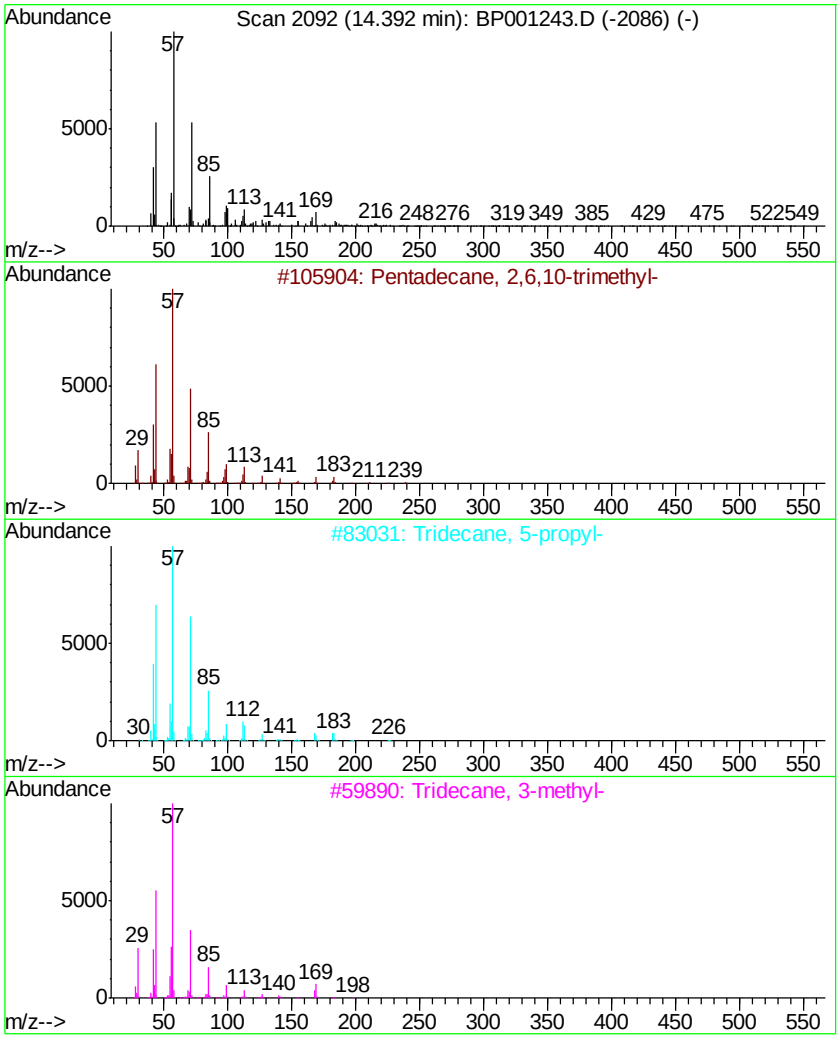
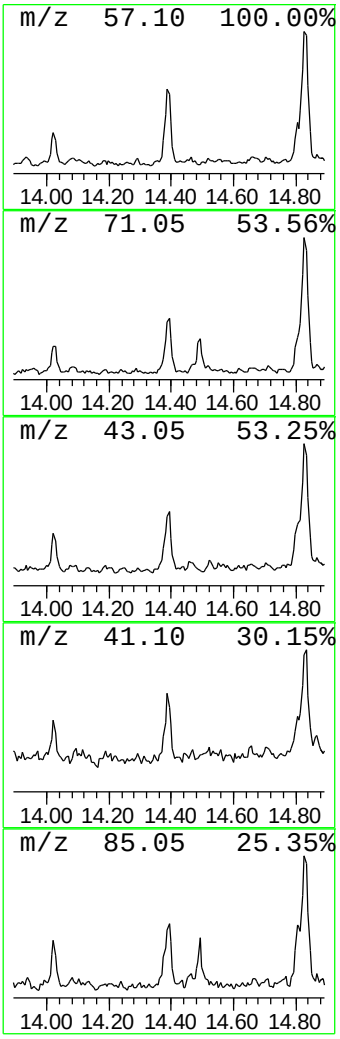
****

Peak Number 5 Pentadecane, 2,6,10-trimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.39	2.24 ng	96491	Acenaphthene-d10	13.20

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	87
2	Tridecane, 5-propyl-	226	C16H34	055045-11-9	83
3	Tridecane, 3-methyl-	198	C14H30	006418-41-3	72
4	Dodecane, 2-methyl-8-propyl-	226	C16H34	055045-07-3	68
5	Decane, 2,6,8-trimethyl-	184	C13H28	062108-26-3	68





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Instrument : BNA_P
ClientSampleId : COMP-36

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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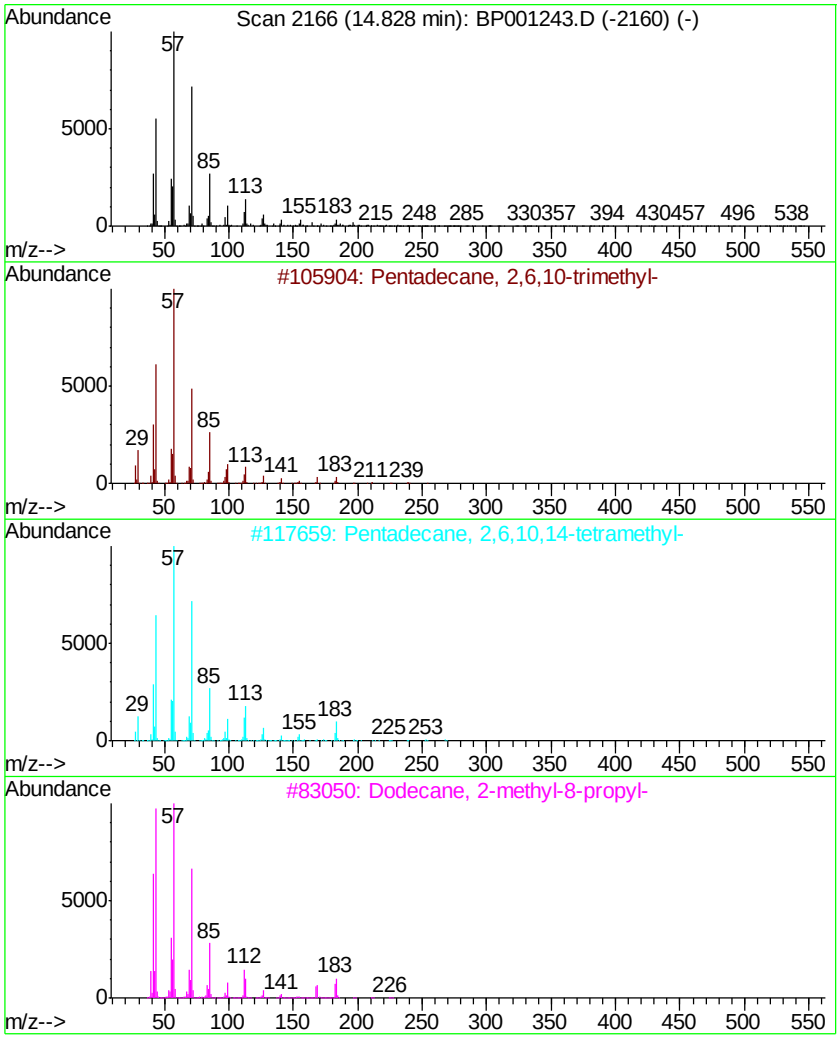
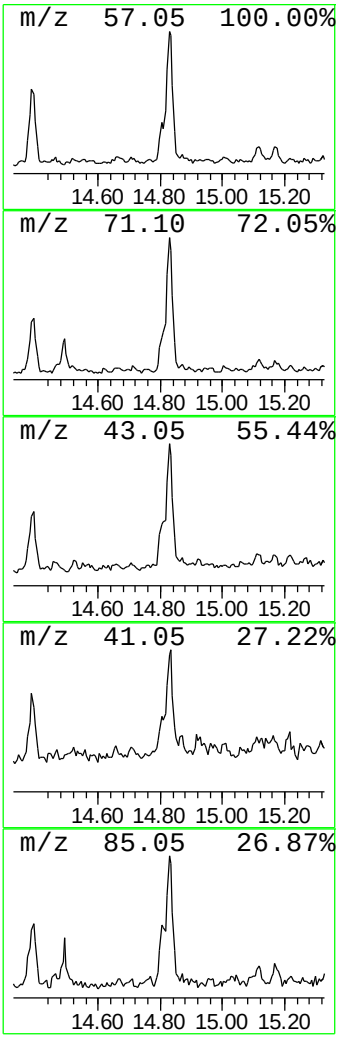
****

Peak Number 6 Pentadecane, 2,6,10,14-tetr... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.83	2.97 ng	121882	Phenanthrene-d10	15.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	94
2	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	91
3	Dodecane, 2-methyl-8-propyl-	226	C16H34	055045-07-3	86
4	Undecane, 6-ethyl-	184	C13H28	017312-60-6	81
5	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	80





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Instrument : BNA_P
ClientSampleId : COMP-36

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

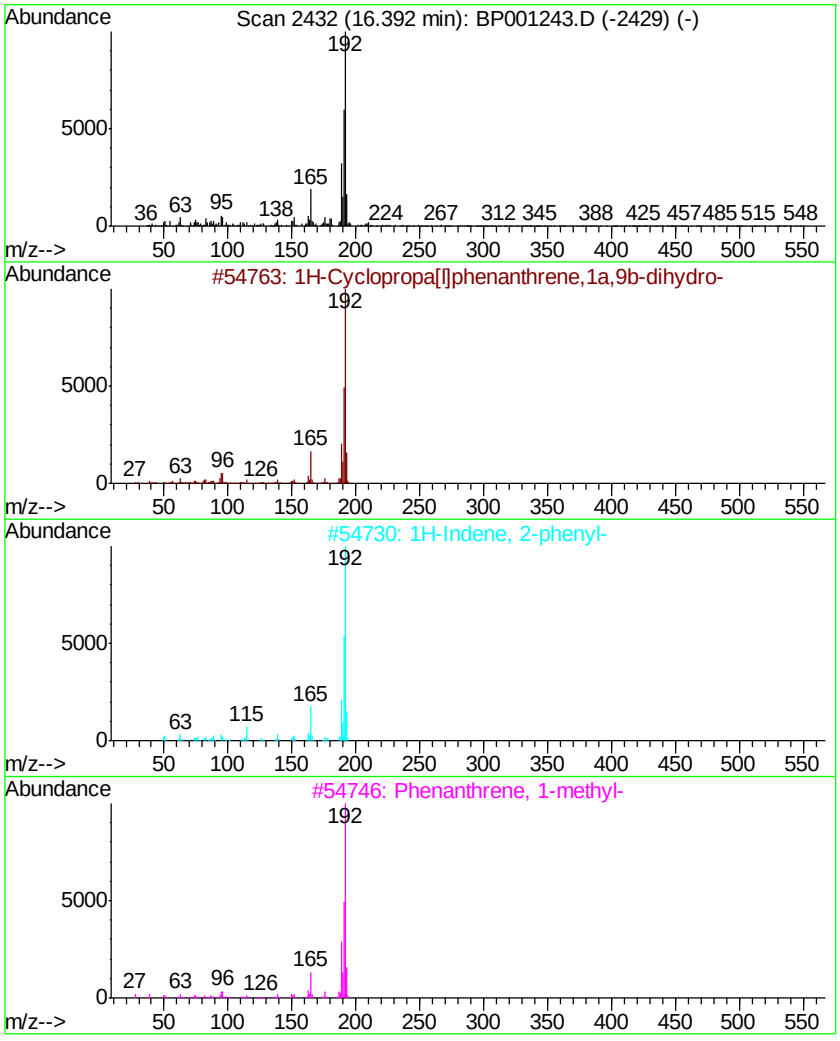
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

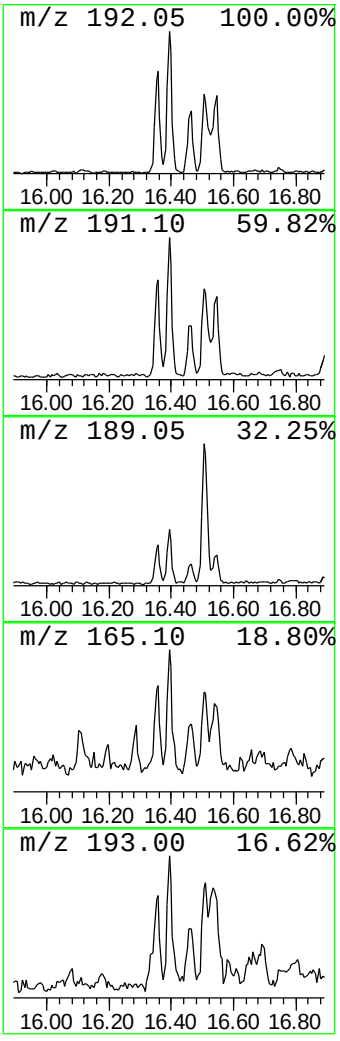
****

Peak Number 7 1H-Cyclopropa[l]phenanthren... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.39	2.70 ng	110567	Phenanthrene-d10	15.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	95
2	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	95
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	95
5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120619\
Data File : BP001243.D
Acq On : 06 Dec 2019 21:51
Operator : JU
Sample : K6135-06 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-36

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

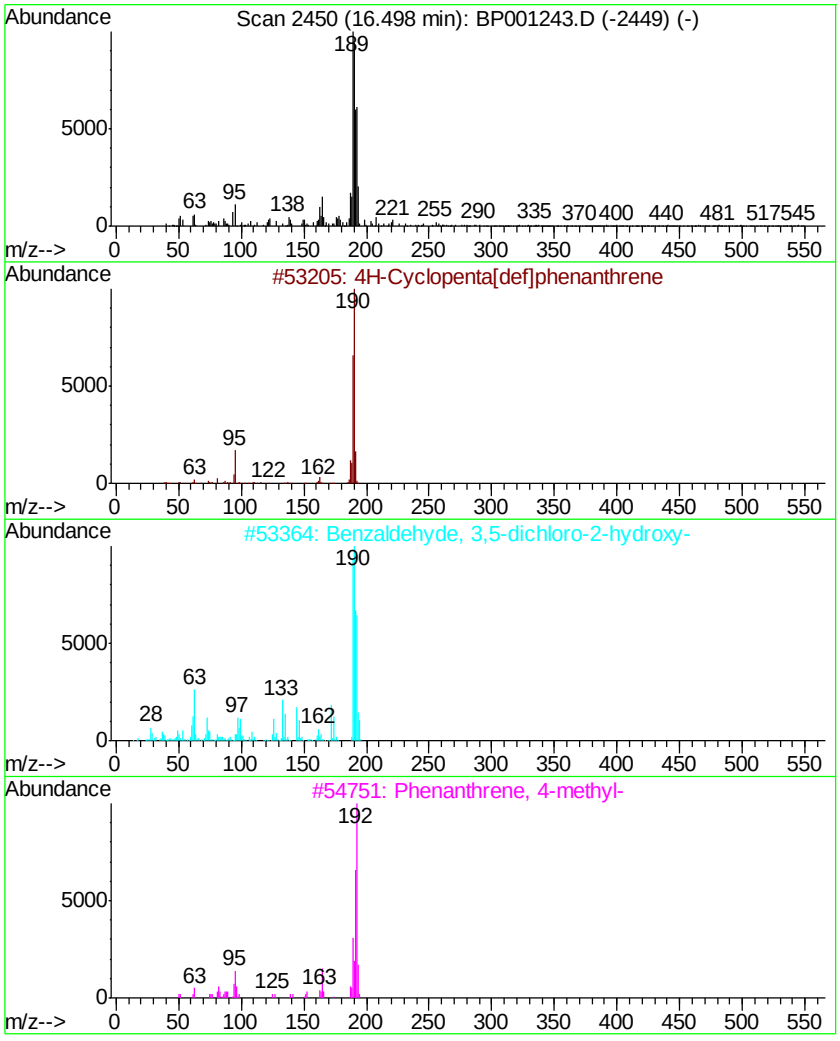
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

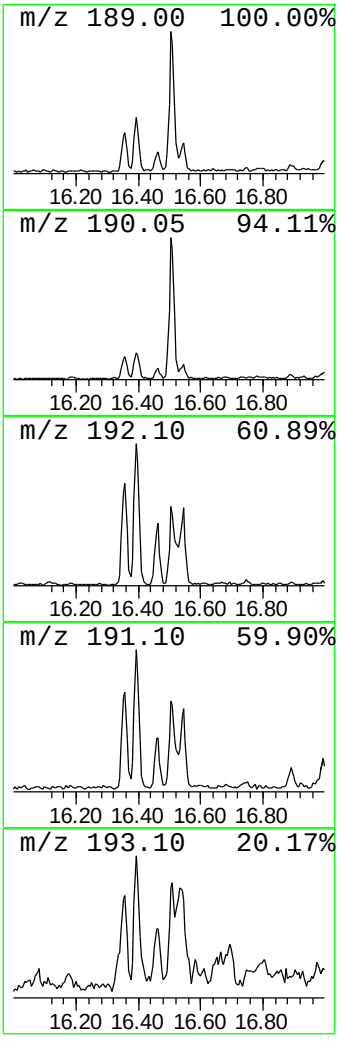
****

Peak Number 8 4H-Cyclopenta[def]phenanthrene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.50	2.08 ng	85298	Phenanthrene-d10	15.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	70
2	Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	50
3	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	42
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	38
5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	38





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120619\
Data File : BP001243.D
Acq On : 06 Dec 2019 21:51
Operator : JU
Sample : K6135-06 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-36

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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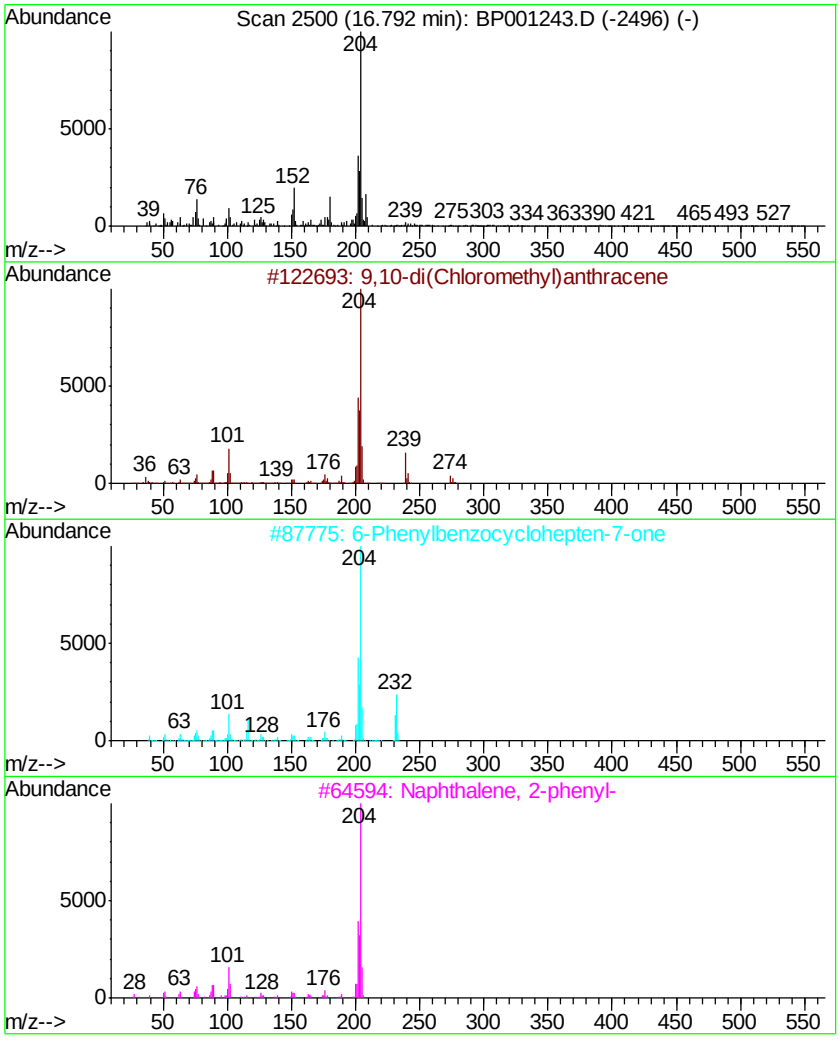
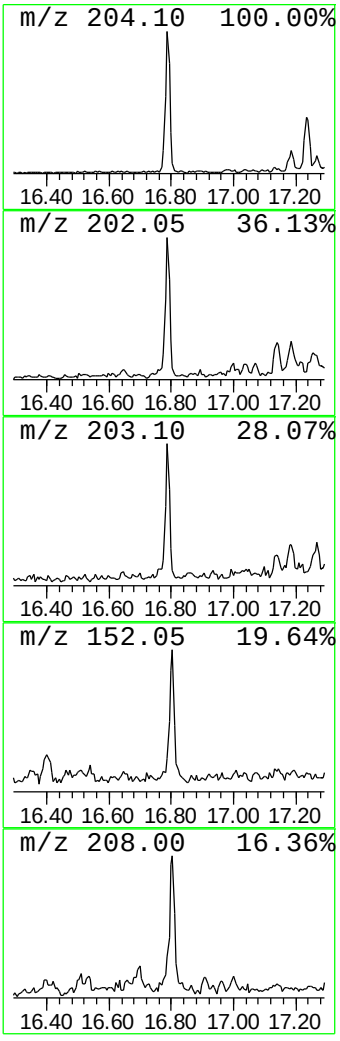
Peak Number 9 9,10-di(Chloromethyl)anthra... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.79	2.49 ng	102157	Phenanthrene-d10	15.60

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-di(Chloromethyl)anthracene	274	C16H12Cl2	010387-13-0	86
2	6-Phenylbenzocyclohepten-7-one	232	C17H12O	093327-56-1	86
3	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	83
4	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	83
5	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	80





Library Search Compound Report

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ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-36

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.M
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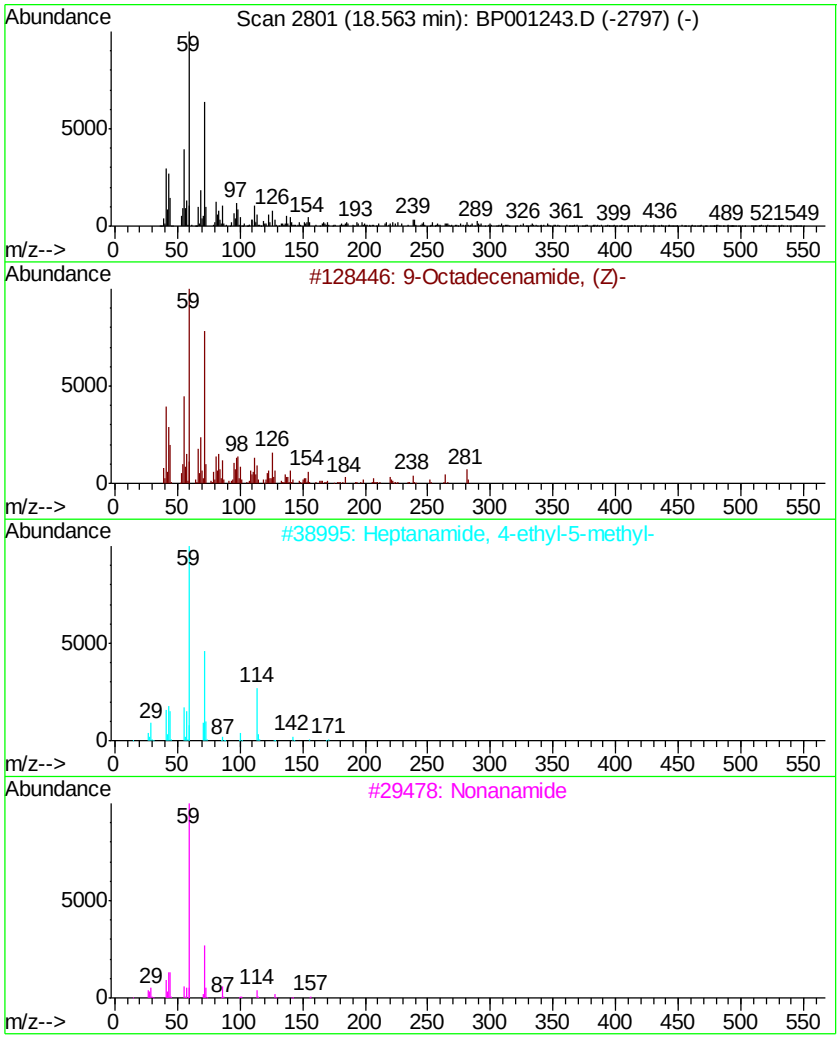
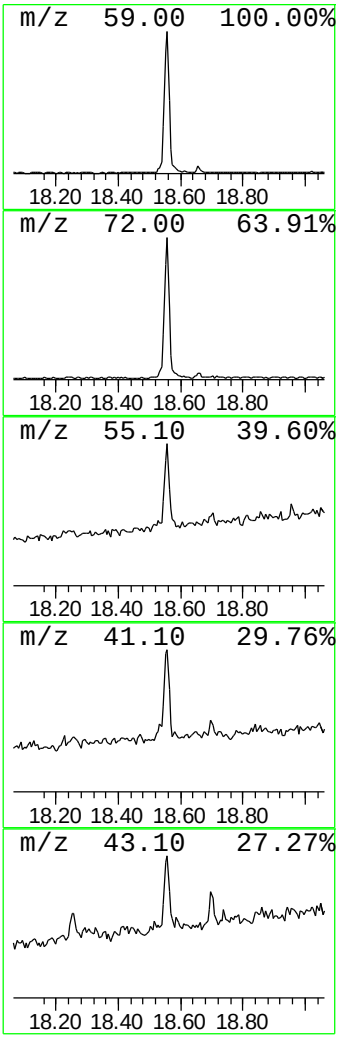
Peak Number 10 9-Octadecenamide, (Z)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.56	3.62 ng	206865	Chrysene-d12	19.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	97
2	Heptanamide, 4-ethyl-5-methyl-	171	C10H21NO	054789-40-1	72
3	Nonanamide	157	C9H19NO	001120-07-6	59
4	Octanamide	143	C8H17NO	000629-01-6	59
5	Hexadecanamide	255	C16H33NO	000629-54-9	50





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120619\
Data File : BP001243.D
Acq On : 06 Dec 2019 21:51
Operator : JU
Sample : K6135-06 2X
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : COMP-36

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP112719.M
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TIC Library : C:\DATABASE\NIST11.L
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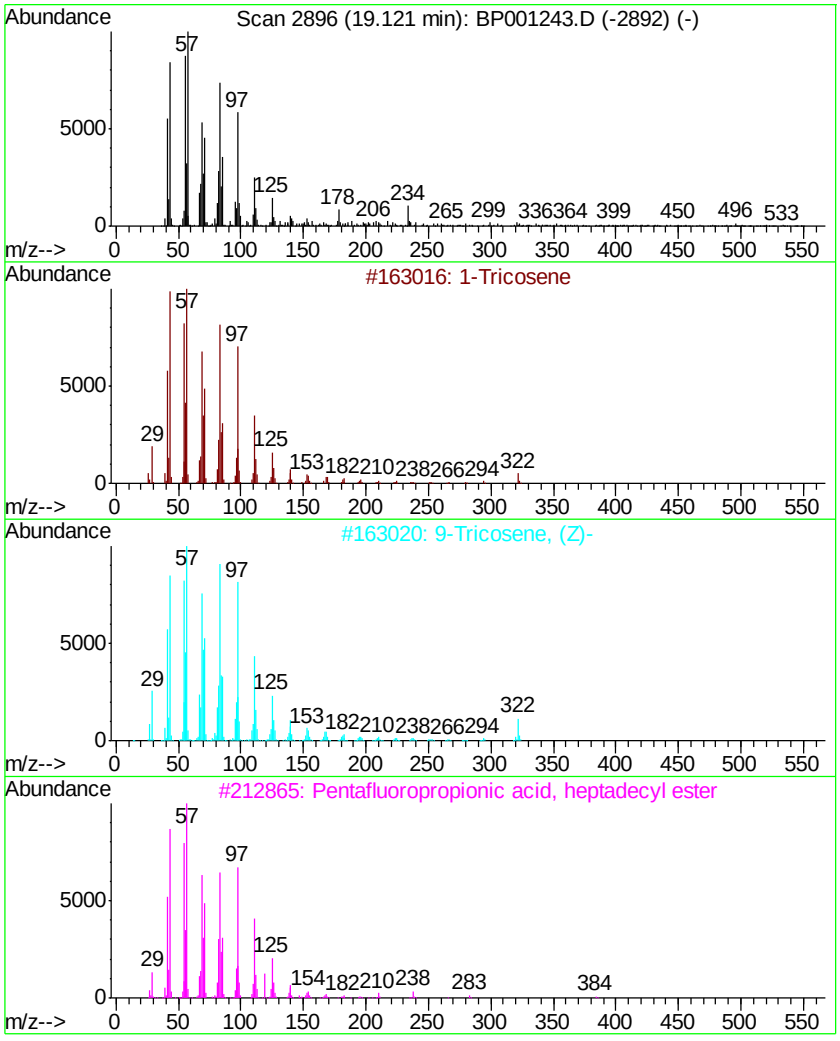
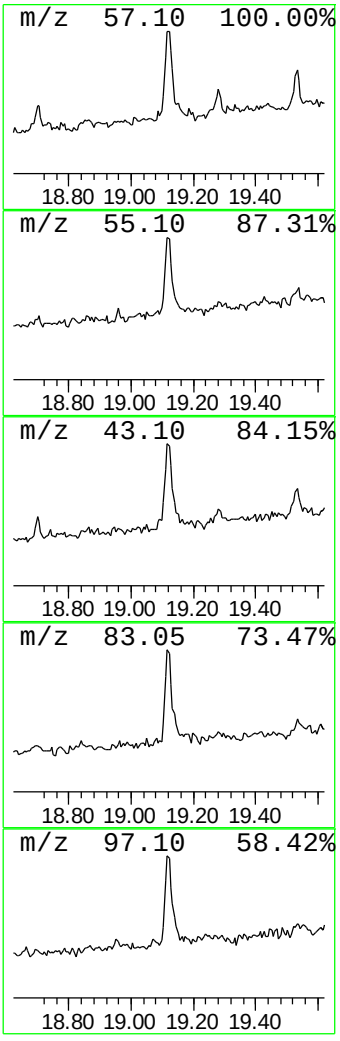
Peak Number 11 1-Tricosene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.12	4.80 ng	273885	Chrysene-d12	19.24

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Tricosene	322	C23H46	018835-32-0	94
2	9-Tricosene, (Z)-	322	C23H46	027519-02-4	94
3	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	93
4	Heptadecyl trifluoroacetate	352	C19H35F3O2	1000351-87-0	93
5	Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP120619\
Data File : BP001243.D
Acq On : 06 Dec 2019 21:51
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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	220078	1	8.31	562071	20.0
unknown7.95	7.95	42.0	ng	1181280	1	8.31	562071	20.0
(+)-2-Bornanone	9.95	3.1	ng	109256	2	10.34	696207	20.0
Pentadecane, 2,6,...	14.39	2.2	ng	96491	3	13.20	860446	20.0
Pentadecane, 2,6,...	14.83	3.0	ng	121882	4	15.60	820437	20.0
1H-Cyclopropa[l]p...	16.39	2.7	ng	110567	4	15.60	820437	20.0
4H-Cyclopenta[def...	16.50	2.1	ng	85298	4	15.60	820437	20.0
9,10-di(Chloromet...	16.79	2.5	ng	102157	4	15.60	820437	20.0
9-Octadecenamide,...	18.56	3.6	ng	206865	5	19.24	1142360	20.0
1-Tricosene	19.12	4.8	ng	273885	5	19.24	1142360	20.0