LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122419\
Data File : BP001404.D
Acq On : 25 Dec 2019 00:04
Operator : JU
Sample : K6413-06
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : K005-AA-WC-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:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP121919.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	4.917	475	481	488	rBV2	14379	22493	1.62%	0.155%
2	5.599	588	597	610	rBV	146587	235912	17.03%	1.627%
3	6.205	694	700	713	rBV	917078	1150862	83.10%	7.937%
4	7.752	956	963	976	rBV	977227	1296118	93.58%	8.938%
5	7.928	987	993	1006	rVB	1083145	1295617	93.55%	8.935%
6	8.281	1048	1053	1062	rBV	285155	339789	24.53%	2.343%
7	8.528	1088	1095	1099	rBV	819909	980405	70.79%	6.761%
8	9.169	1197	1204	1208	rBV	646294	804967	58.12%	5.551%
9	10.316	1394	1399	1404	rBV	328253	402961	29.10%	2.779%
10	12.099	1693	1702	1707	rBV	1166190	1384977	100.00%	9.551%
11	12.763	1811	1815	1820	rBV	62086	73349	5.30%	0.506%
12	12.828	1821	1826	1839	rVB6	7691	17526	1.27%	0.121%
13	13.181	1879	1886	1891	rBV	379693	453155	32.72%	3.125%
14	13.228	1891	1894	1899	rVB	15357	19465	1.41%	0.134%
15	14.075	2035	2038	2046	rVB	13672	18748	1.35%	0.129%
16	14.363	2079	2087	2094	rBV8	8321	18040	1.30%	0.124%
17	14.475	2098	2106	2117	rBV	731848	910481	65.74%	6.279%
18	14.810	2160	2163	2170	rVB5	10777	17193	1.24%	0.119%
19	15.581	2289	2294	2297	rBV	356020	415509	30.00%	2.865%
20	15.616	2297	2300	2305	rVB	254344	264618	19.11%	1.825%
21	15.692	2309	2313	2319	rVB	60788	70135	5.06%	0.484%
22	16.334	2418	2422	2426	rBV	50642	60844	4.39%	0.420%
23	16.375	2426	2429	2433	rVB	61889	65676	4.74%	0.453%
24	16.440	2437	2440	2442	rBV	19282	23196	1.67%	0.160%
25	16.475	2442	2446	2452	rVV2	104083	213863	15.44%	1.475%
26	16.522	2452	2454	2459	rVB	36785	42235	3.05%	0.291%
27	16.769	2492	2496	2505	rVV2	39202	59228	4.28%	0.408%
28	16.975	2528	2531	2536	rBV2	13105	15096	1.09%	0.104%
29	17.122	2552	2556	2560	rBV2	28341	42075	3.04%	0.290%
30	17.163	2560	2563	2567	rVV2	16519	21167	1.53%	0.146%
31	17.328	2586	2591	2595	rBV	263709	282750	20.42%	1.950%
32	17.557	2627	2630	2635	rBV2	36529	45335	3.27%	0.313%
33	17.634	2638	2643	2648	rVB	341176	391481	28.27%	2.700%
34	17.863	2677	2682	2686	rBV	1059006	1178242	85.07%	8.126%

LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122419\
Data File : BP001404.D
Acq On : 25 Dec 2019 00:04
Operator : JU
Sample : K6413-06
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : K005-AA-WC-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:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP121919.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

35	17.945	2692	2696	2700	rBV2	18552	28220	2.04%	0.195%
36	18.086	2718	2720	2727	rVB	36788	40229	2.90%	0.277%
37	18.175	2732	2735	2739	rVB2	19680	23271	1.68%	0.160%
38	18.222	2740	2743	2748	rVV2	41543	59758	4.31%	0.412%
39	18.339	2760	2763	2767	rBV	25104	27226	1.97%	0.188%
40	18.381	2767	2770	2775	rVV	26844	33238	2.40%	0.229%
41	18.681	2819	2821	2824	rVB	24215	19963	1.44%	0.138%
42	18.933	2862	2864	2867	rBV	22234	26423	1.91%	0.182%
43	19.104	2889	2893	2900	rBV2	110112	173870	12.55%	1.199%
44	19.222	2907	2913	2917	rBV2	357867	544012	39.28%	3.752%
45	19.257	2917	2919	2923	rVB	126960	114725	8.28%	0.791%
46	19.516	2960	2963	2966	rVB	33726	36895	2.66%	0.254%
47	20.280	3090	3093	3099	rVB2	47397	56124	4.05%	0.387%
48	20.510	3128	3132	3135	rBV	82931	129367	9.34%	0.892%
49	20.669	3157	3159	3164	rVB3	43759	47071	3.40%	0.325%
50	20.851	3187	3190	3197	rVB	62745	84126	6.07%	0.580%
51	20.922	3198	3202	3206	rVB	85110	110899	8.01%	0.765%
52	20.992	3210	3214	3219	rBV	243719	311537	22.49%	2.148%

Sum of corrected areas: 14500462

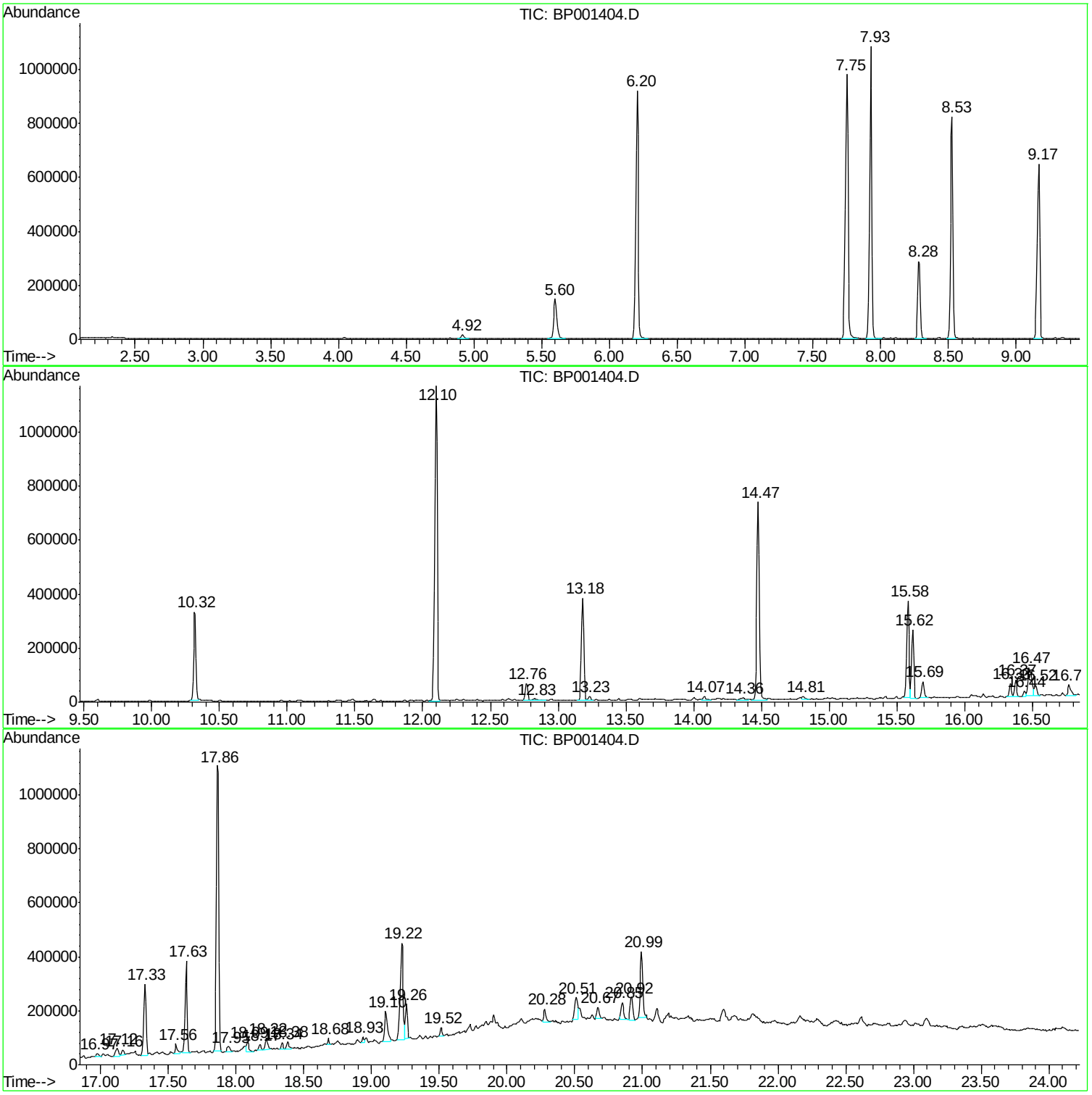
LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122419\
Data File : BP001404.D
Acq On : 25 Dec 2019 00:04
Operator : JU
Sample : K6413-06
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : K005-AA-WC-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP121919.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 P\DATA\BP122419\
Data File : BP001404.D
Acq On : 25 Dec 2019 00:04
Operator : JU
Sample : K6413-06
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : K005-AA-WC-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP121919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

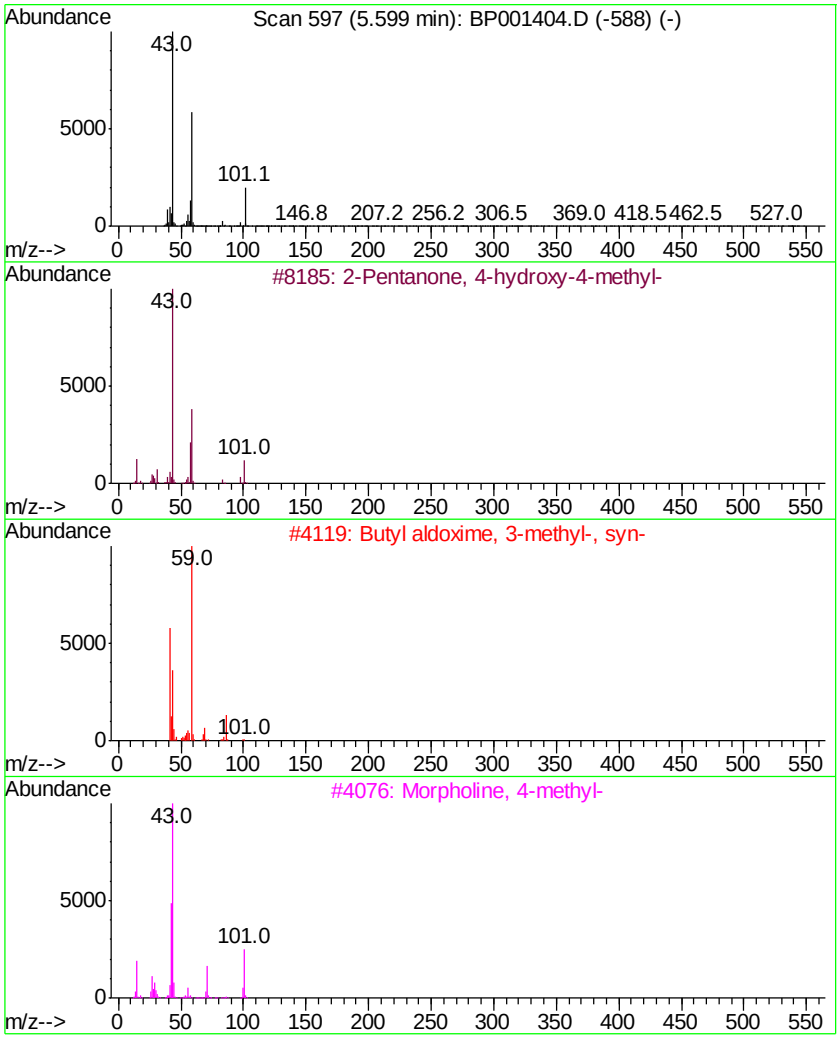
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

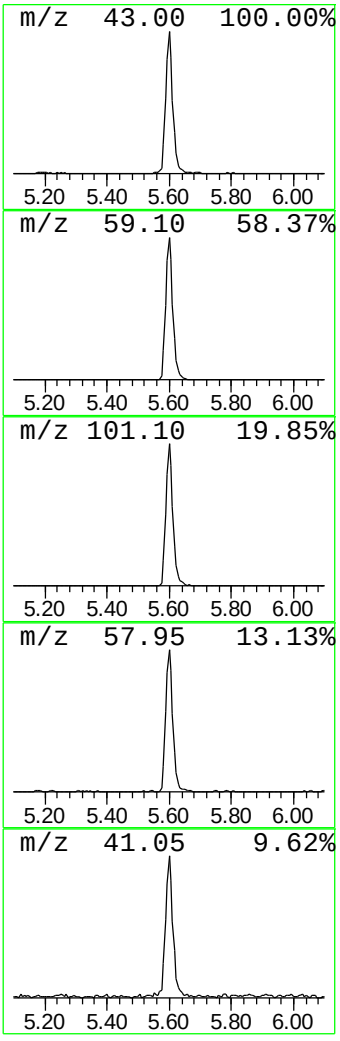
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.60	13.89 ng	235912	1,4-Dichlorobenzene-d4	8.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2	Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	17
3	Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
4	Butane	58	C4H10	000106-97-8	9
5	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122419\
Data File : BP001404.D
Acq On : 25 Dec 2019 00:04
Operator : JU
Sample : K6413-06
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : K005-AA-WC-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP121919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

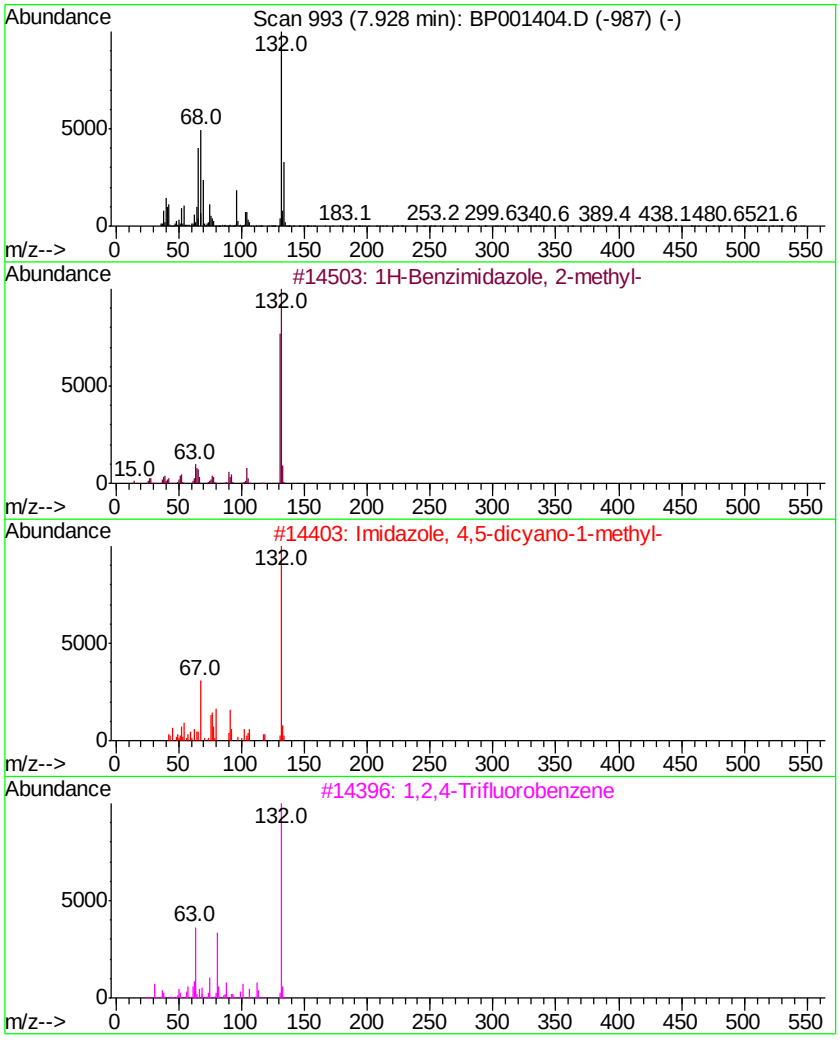
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

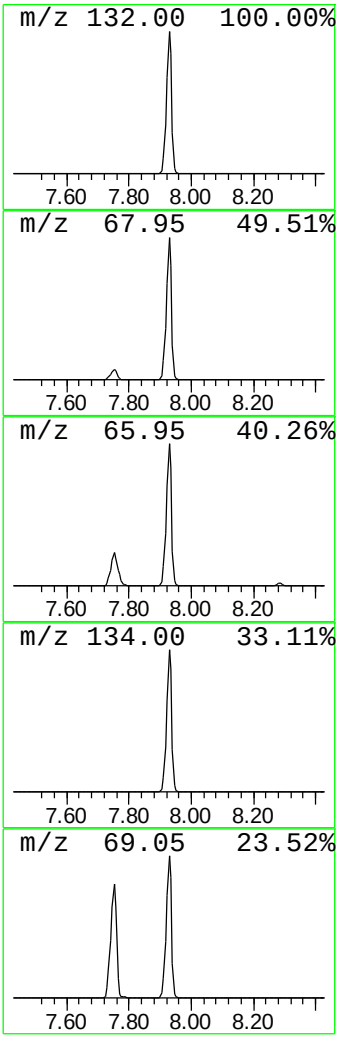
****

Peak Number 2 unknown7.93 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.93	76.26 ng	1295620	1,4-Dichlorobenzene-d4	8.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	14
2	Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	10
3	1,2,4-Trifluorobenzene	132	C6H3F3	000367-23-7	10
4	1,3-Cyclopentanedione, 2-chloro-	132	C5H5ClO2	014203-19-1	10
5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	10





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122419\
Data File : BP001404.D
Acq On : 25 Dec 2019 00:04
Operator : JU
Sample : K6413-06
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : K005-AA-WC-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP121919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

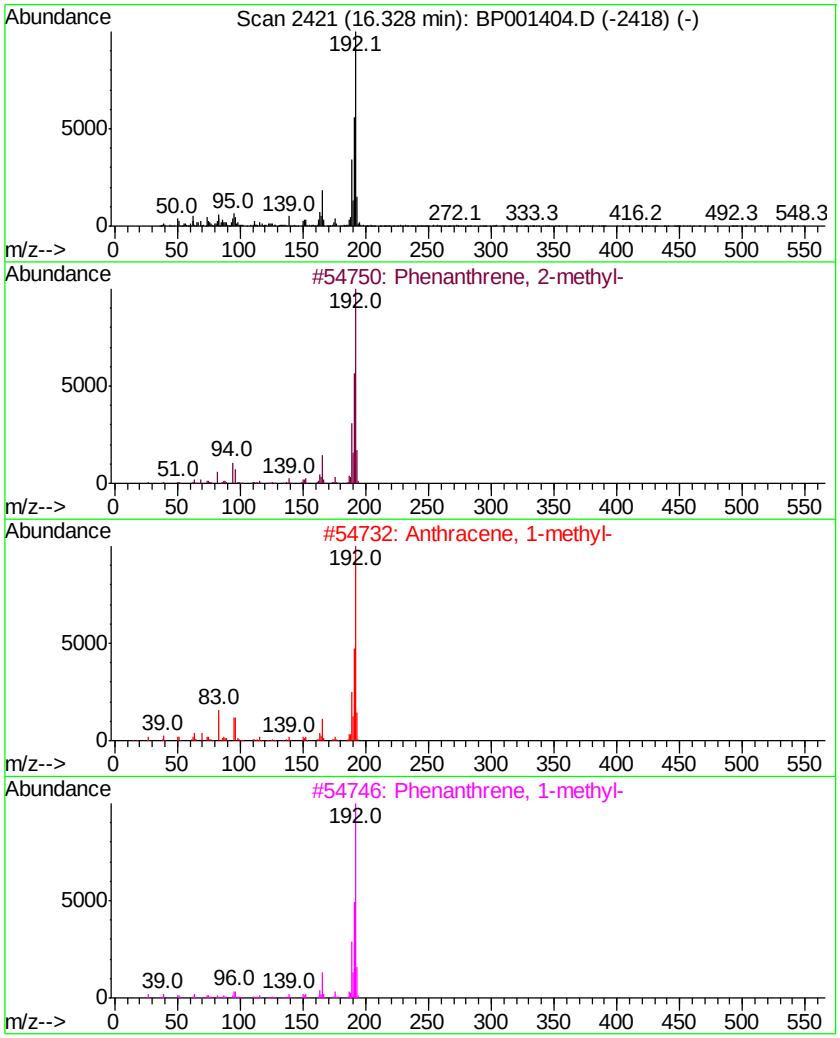
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

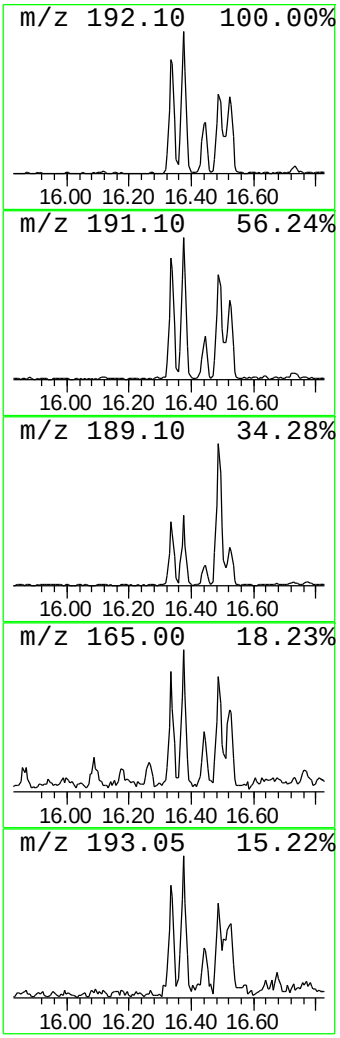
****

Peak Number 4 Phenanthrene, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.33	2.93 ng	60844	Phenanthrene-d10	15.58

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
2	Anthracene, 1-methyl-	192	C15H12	000610-48-0	90
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	90
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	87
5	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	87





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122419\
Data File : BP001404.D
Acq On : 25 Dec 2019 00:04
Operator : JU
Sample : K6413-06
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : K005-AA-WC-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP121919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

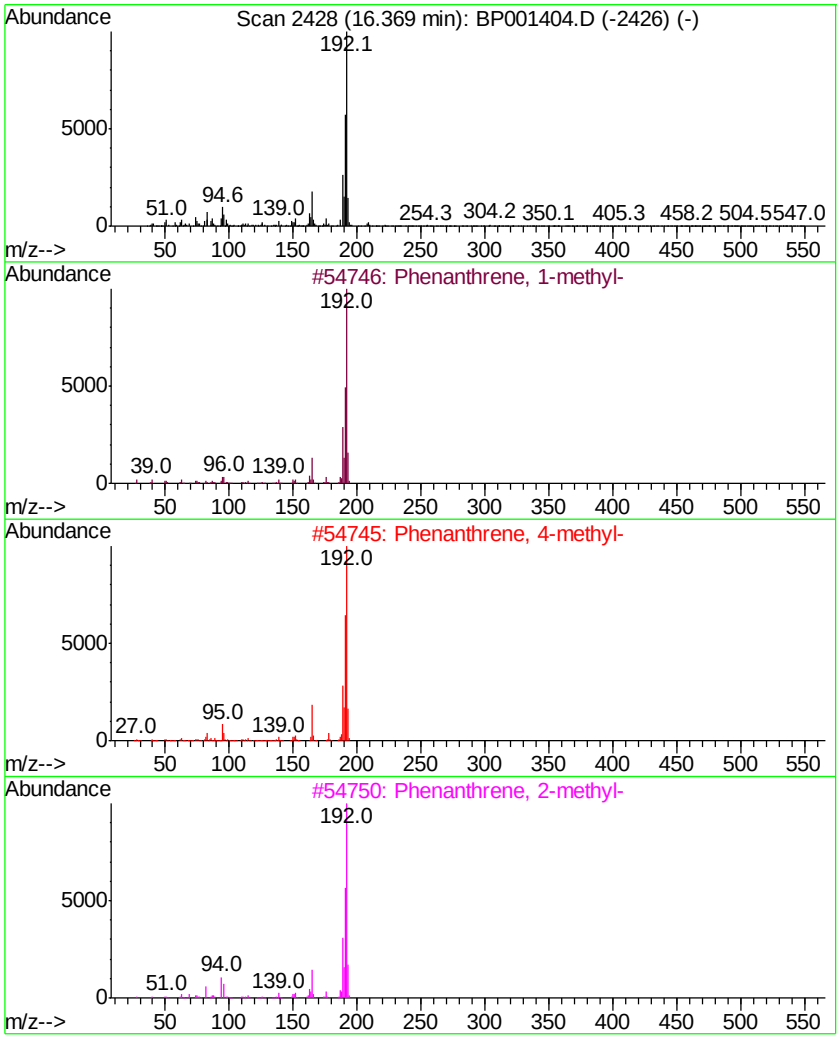
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

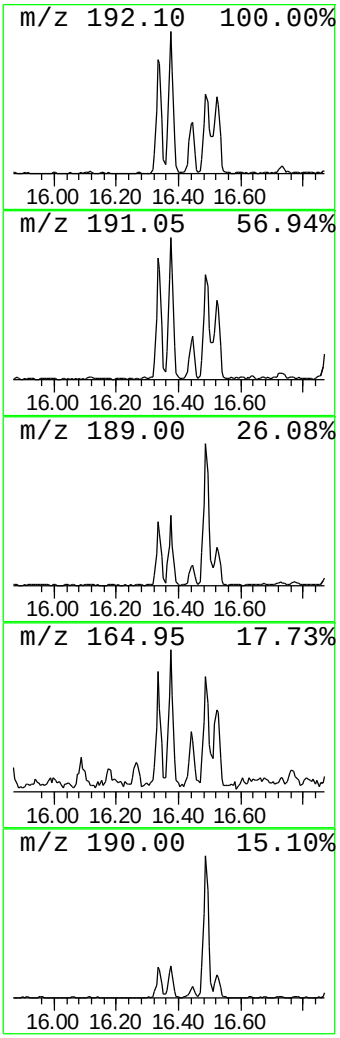
****

Peak Number 5 Phenanthrene, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.37	3.16 ng	65676	Phenanthrene-d10	15.58

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
2	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	96
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
4	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	94
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	91





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122419\
Data File : BP001404.D
Acq On : 25 Dec 2019 00:04
Operator : JU
Sample : K6413-06
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : K005-AA-WC-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP121919.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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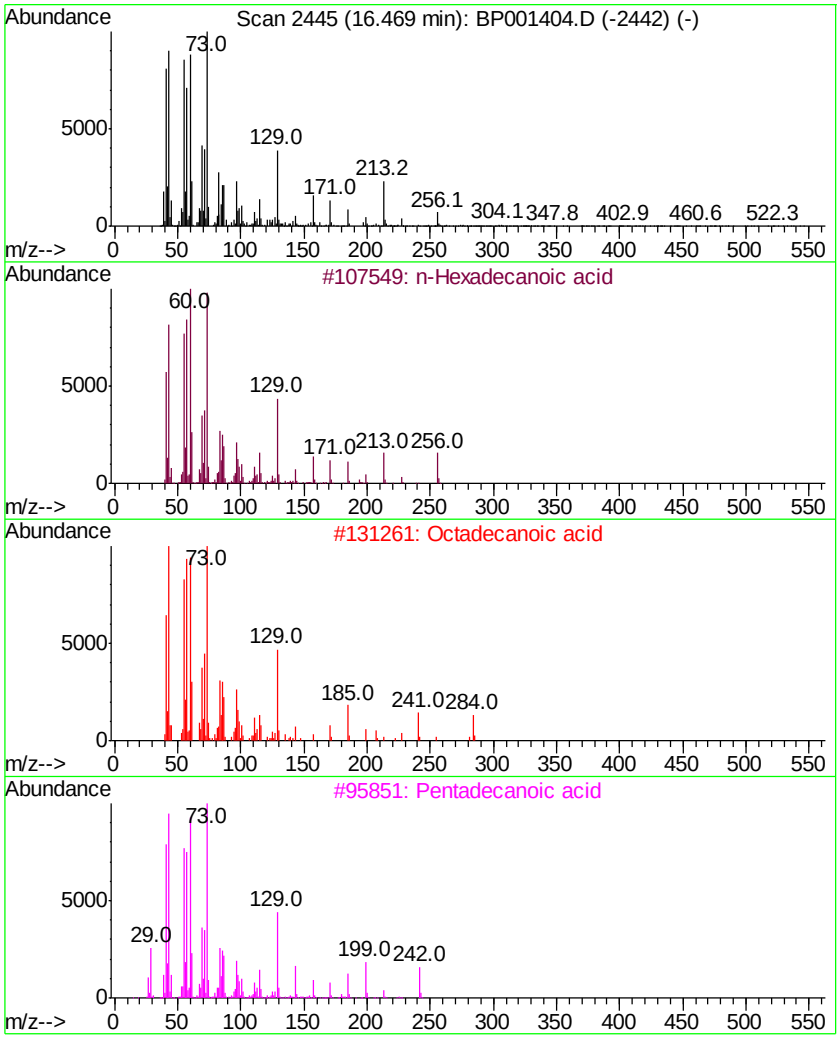
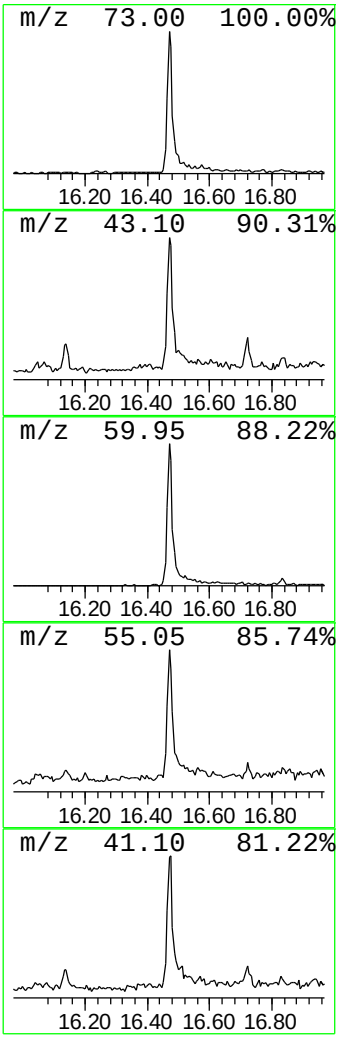
Peak Number 6 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.47	10.29 ng	213863	Phenanthrene-d10	15.58

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Octadecanoic acid	284	C18H36O2	000057-11-4	76
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	68
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	58
5	Tridecanoic acid	214	C13H26O2	000638-53-9	53





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122419\
Data File : BP001404.D
Acq On : 25 Dec 2019 00:04
Operator : JU
Sample : K6413-06
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : K005-AA-WC-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP121919.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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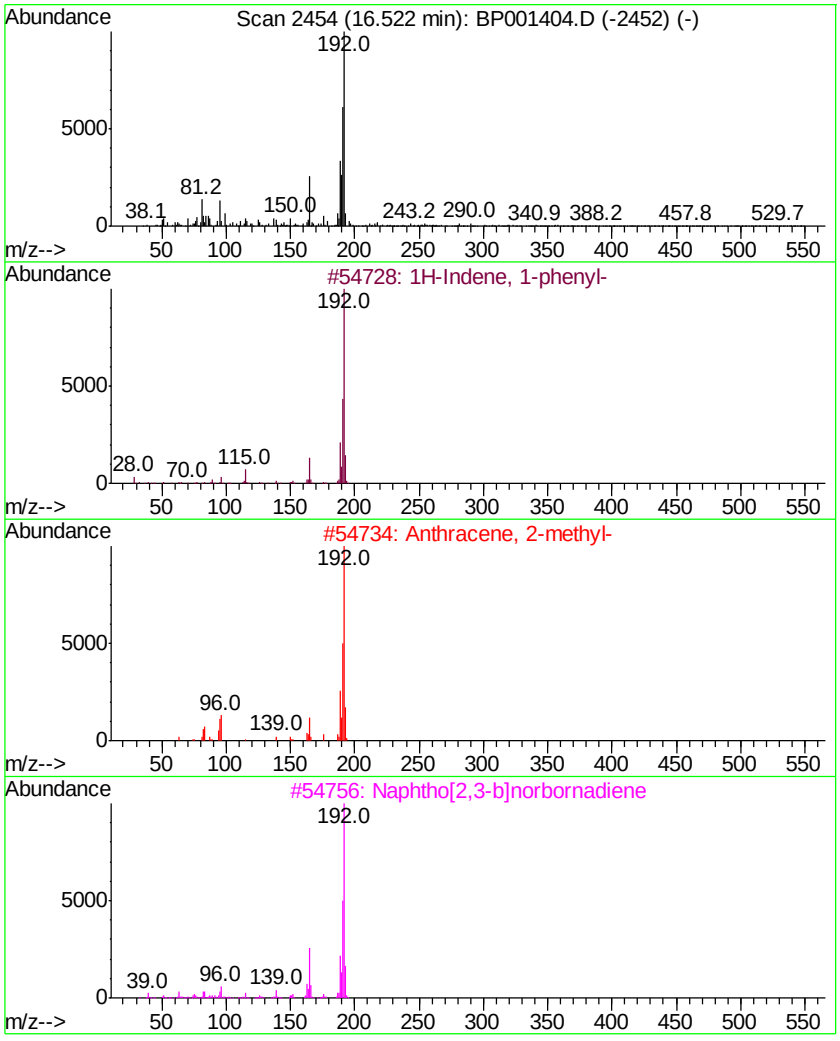
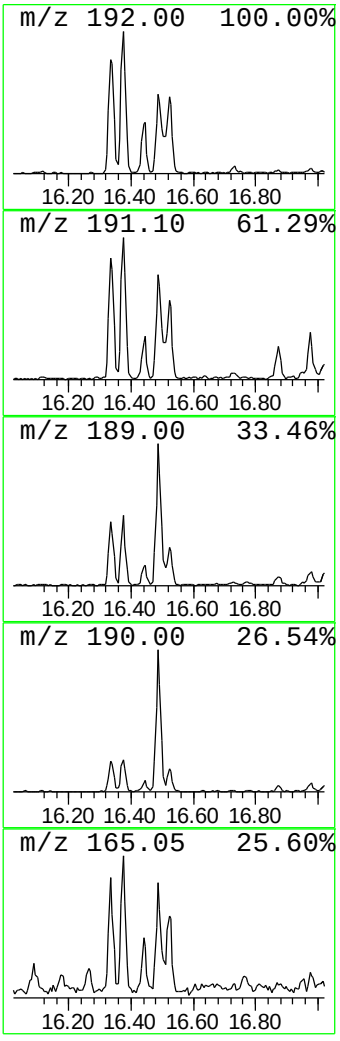
Peak Number 7 1H-Indene, 1-phenyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.52	2.03 ng	42235	Phenanthrene-d10	15.58

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 1-phenyl-	192	C15H12	001961-96-2	64
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	62
3	Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	59
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	59
5	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	59





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122419\
Data File : BP001404.D
Acq On : 25 Dec 2019 00:04
Operator : JU
Sample : K6413-06
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : K005-AA-WC-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP121919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

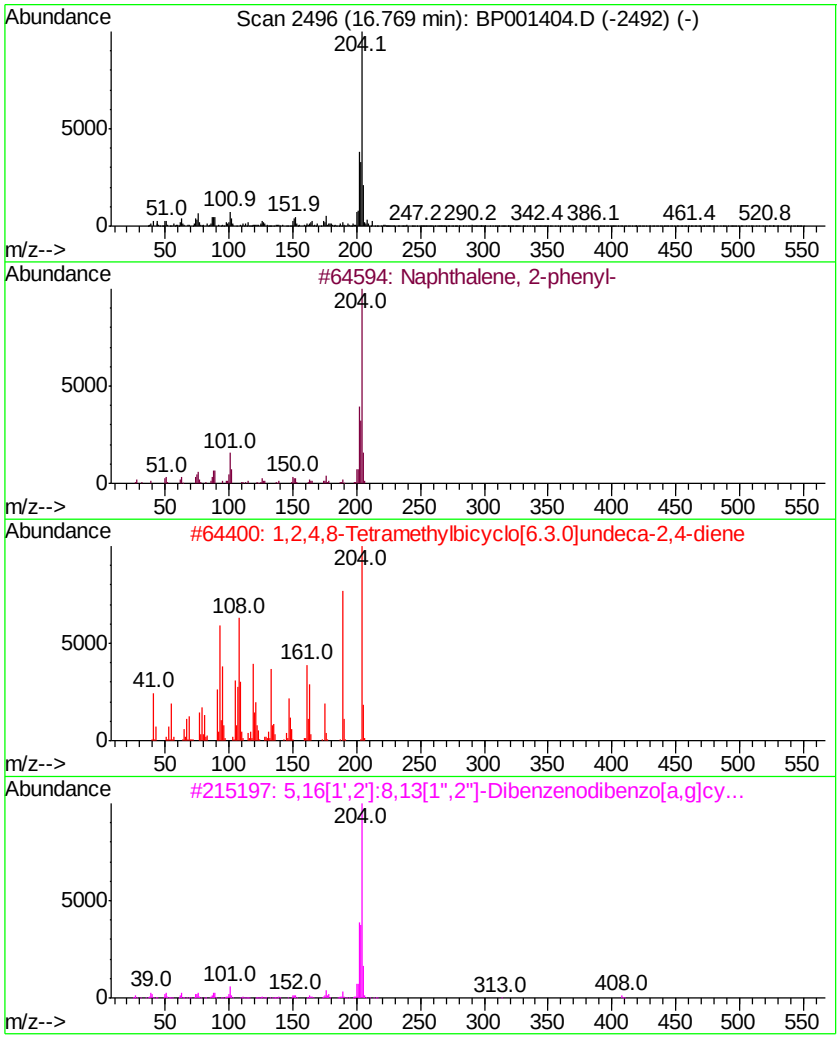
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

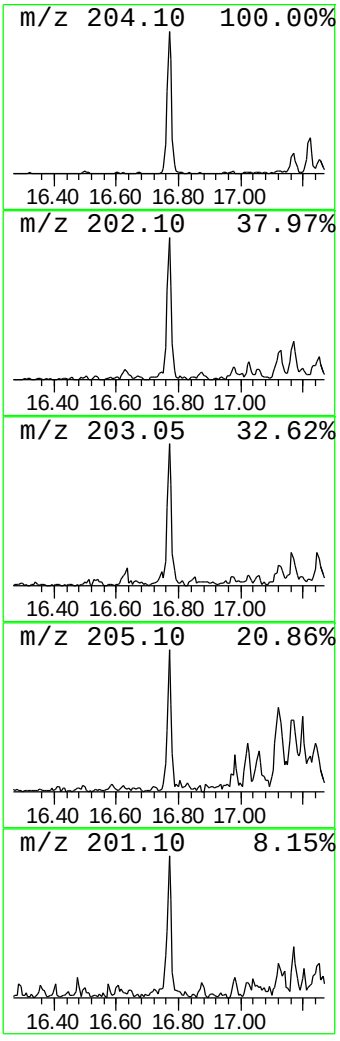
****

Peak Number 8 Naphthalene, 2-phenyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.77	2.85 ng	59228	Phenanthrene-d10	15.58

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	90
2	1,2,4,8-Tetramethylbicyclo[6.3.0...	204	C15H24	137235-51-9	90
3	5,16[1',2']:8,13[1'',2'']-Dibenz...	408	C32H24	005672-97-9	90
4	9,10-Bis(bromomethyl)anthracene	362	C16H12Br2	034373-96-1	83
5	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	72





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122419\
Data File : BP001404.D
Acq On : 25 Dec 2019 00:04
Operator : JU
Sample : K6413-06
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : K005-AA-WC-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP121919.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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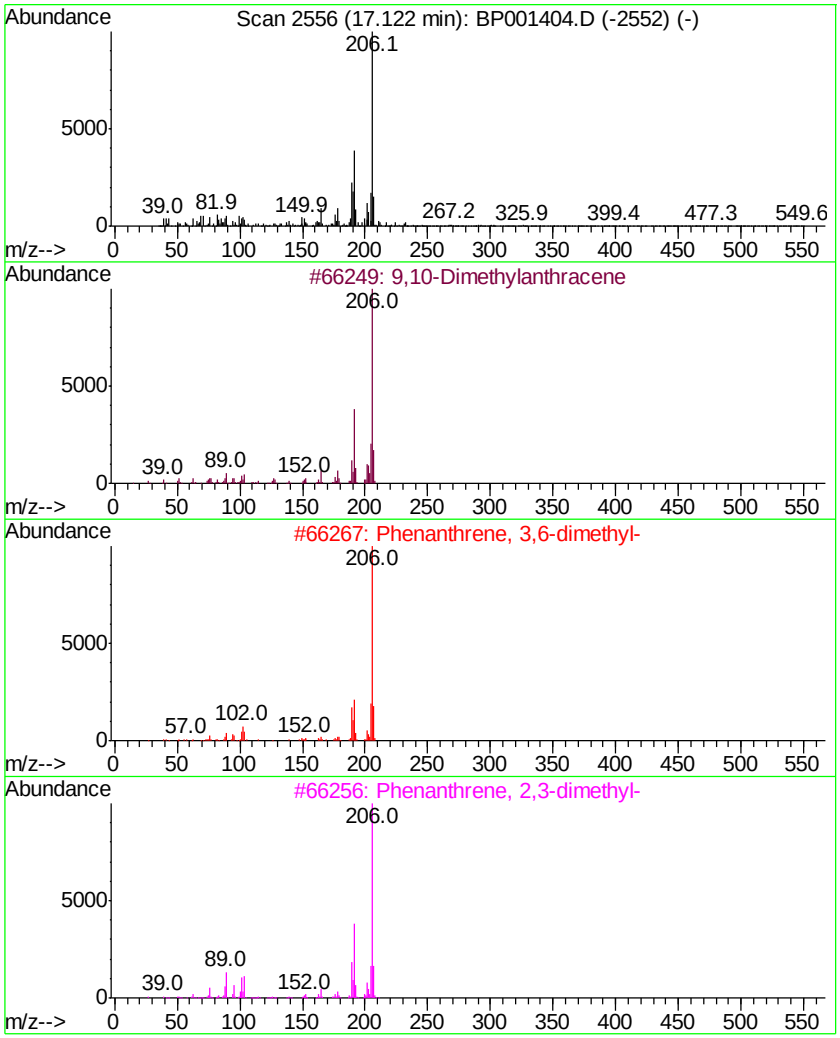
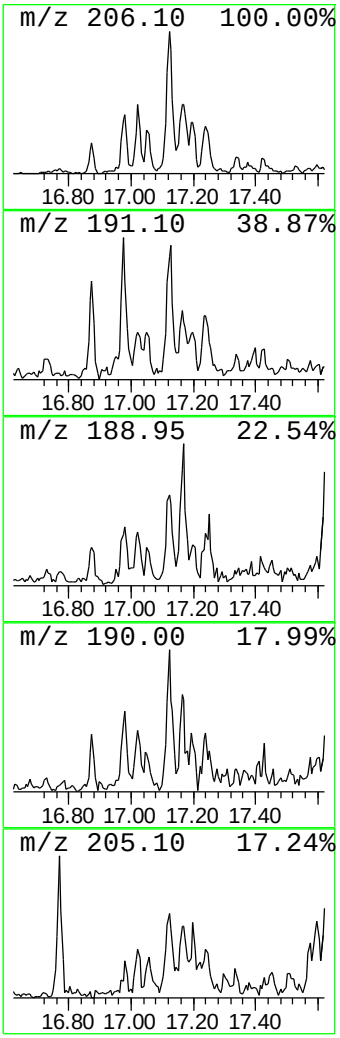
Peak Number 9 9,10-Dimethylanthracene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.12	2.03 ng	42075	Phenanthrene-d10	15.58

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Dimethylanthracene	206	C16H14	000781-43-1	90
2	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	87
3	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	87
4	Naphthalene, 1,2-dihydro-4-phenyl-	206	C16H14	007469-40-1	83
5	Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	83





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Data File : BP001404.D
Acq On : 25 Dec 2019 00:04
Operator : JU
Sample : K6413-06
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : K005-AA-WC-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP121919.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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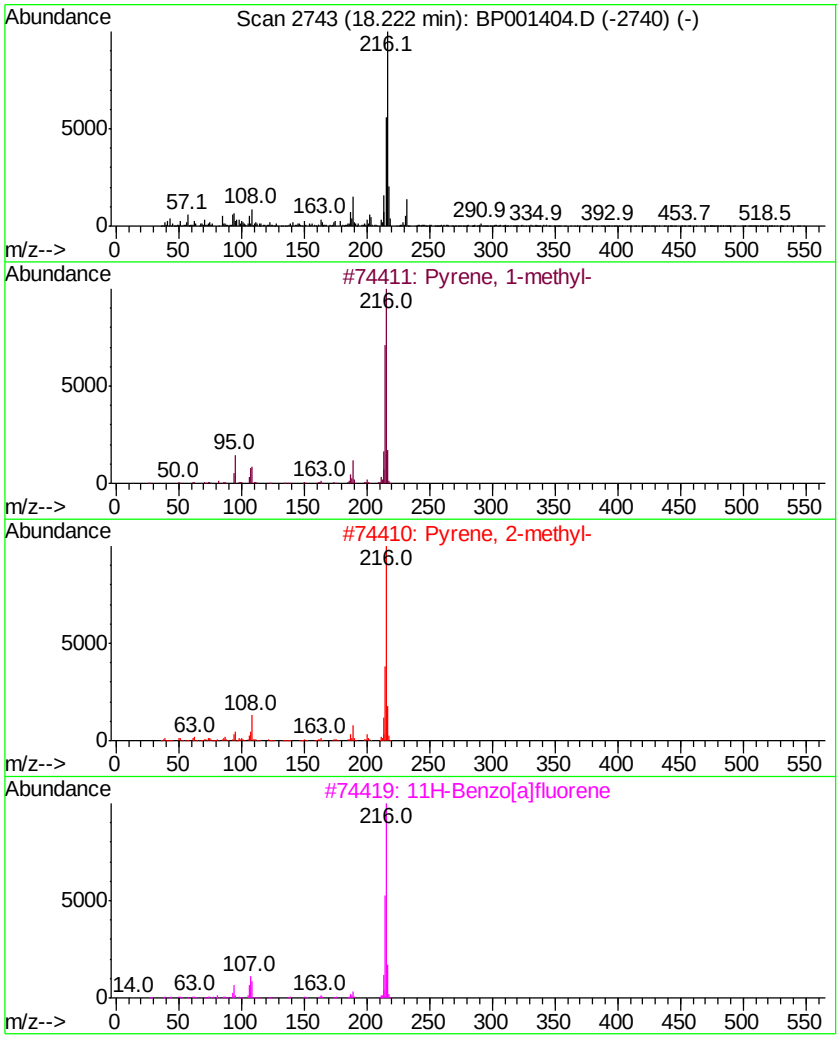
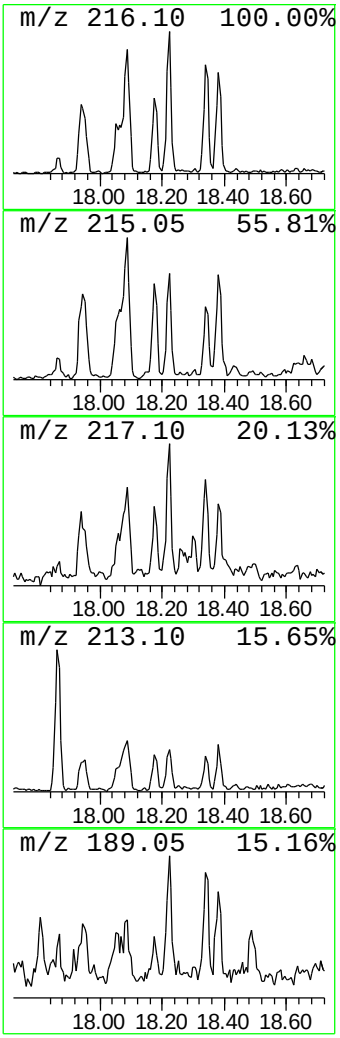
Peak Number 10 Pyrene, 1-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.22	2.20 ng	59758	Chrysene-d12	19.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
2	Pyrene, 2-methyl-	216	C17H12	003442-78-2	93
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	93
4	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	83
5	Pyrene, 4-methyl-	216	C17H12	003353-12-6	81





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Data File : BP001404.D
Acq On : 25 Dec 2019 00:04
Operator : JU
Sample : K6413-06
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : K005-AA-WC-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP121919.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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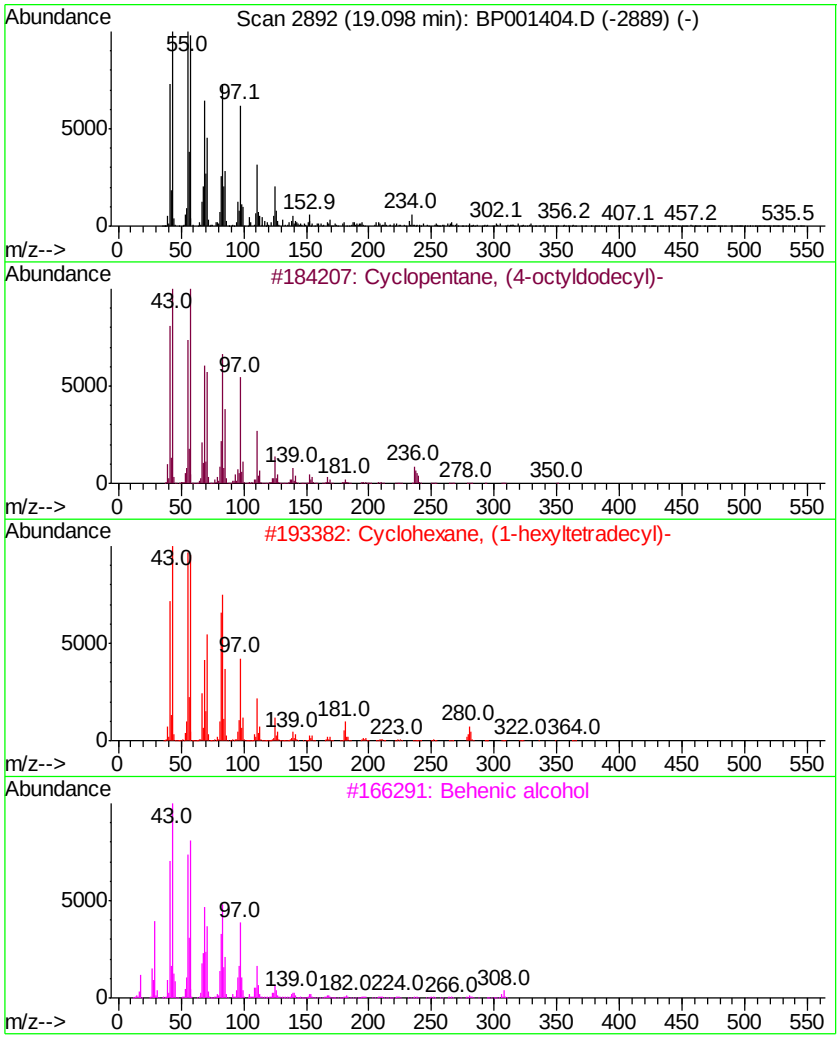
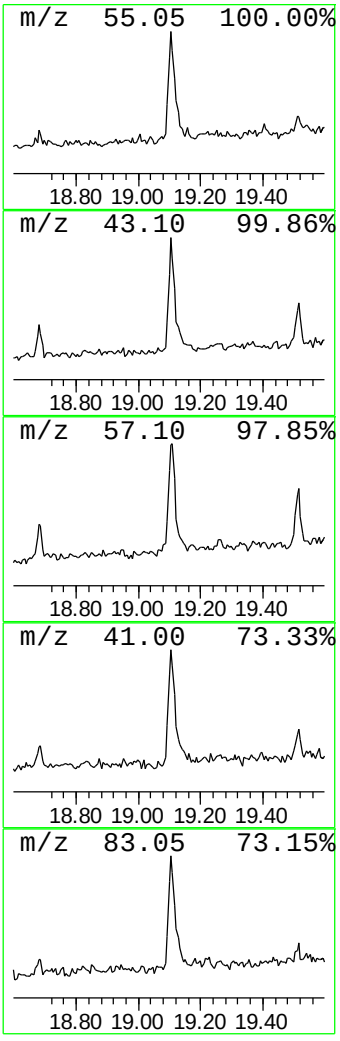
Peak Number 11 Cyclopentane, (4-octyldodec... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.10	6.39 ng	173870	Chrysene-d12	19.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, (4-octyldodecyl)-	350	C25H50	005638-09-5	93
2	Cyclohexane, (1-hexyltetradecyl)-	364	C26H52	004443-60-1	93
3	Behenic alcohol	326	C22H46O	000661-19-8	90
4	17-Pentatriacontene	491	C35H70	006971-40-0	90
5	1-Octadecanethiol	286	C18H38S	002885-00-9	89





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122419\
Data File : BP001404.D
Acq On : 25 Dec 2019 00:04
Operator : JU
Sample : K6413-06
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : K005-AA-WC-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP121919.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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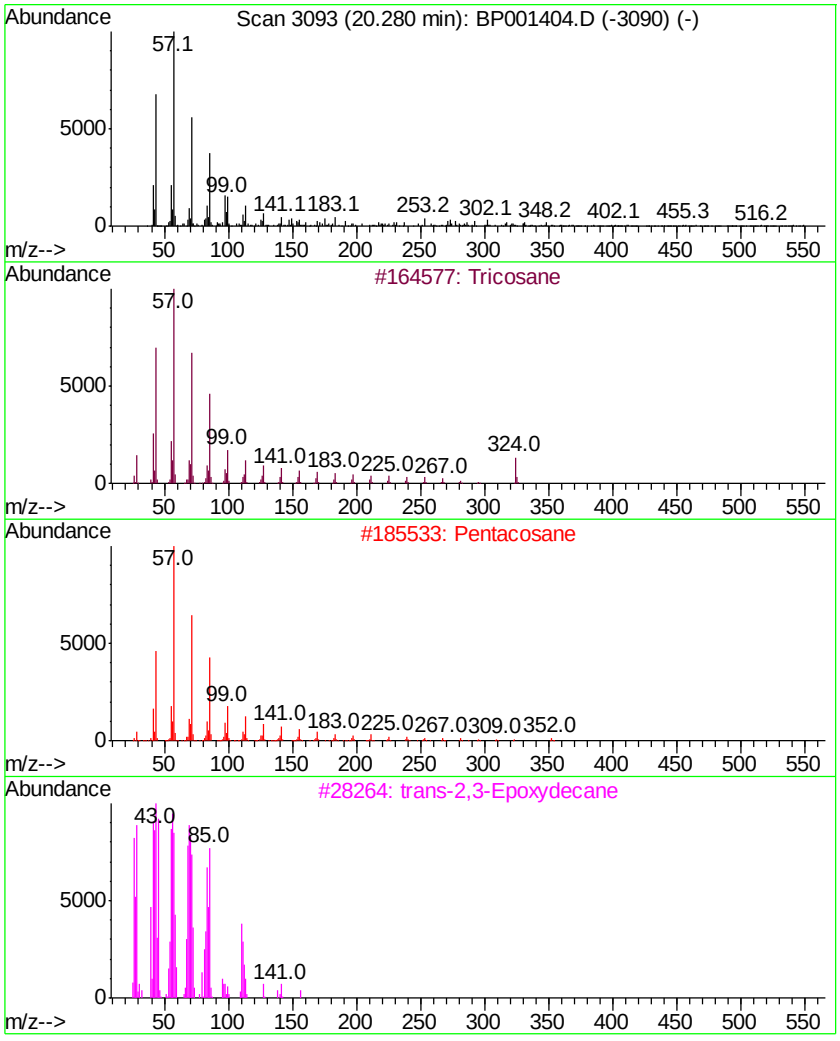
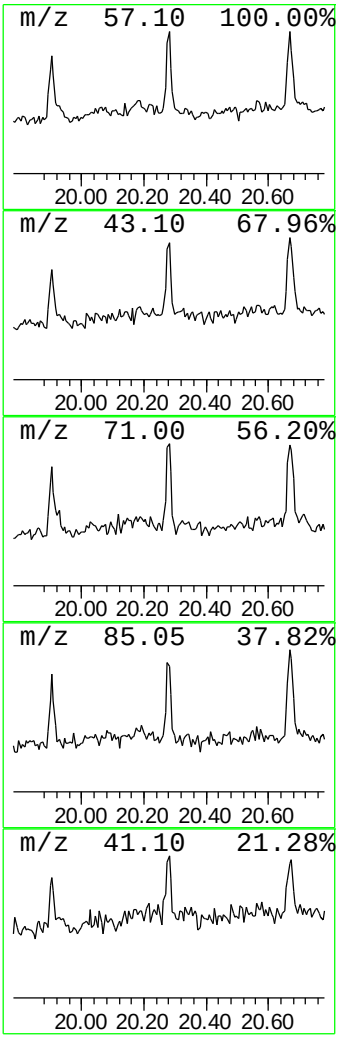
Peak Number 12 Tricosane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.28	3.60 ng	56124	Perylene-d12	20.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tricosane	324	C23H48	000638-67-5	91
2	Pentacosane	352	C25H52	000629-99-2	91
3	trans-2,3-Epoxydecane	156	C10H20O	054125-39-2	90
4	Nonadecane, 1-chloro-	302	C19H39Cl	062016-76-6	87
5	Eicosane	282	C20H42	000112-95-8	83





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122419\
Data File : BP001404.D
Acq On : 25 Dec 2019 00:04
Operator : JU
Sample : K6413-06
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : K005-AA-WC-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP121919.M
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TIC Library : C:\DATABASE\NIST11.L
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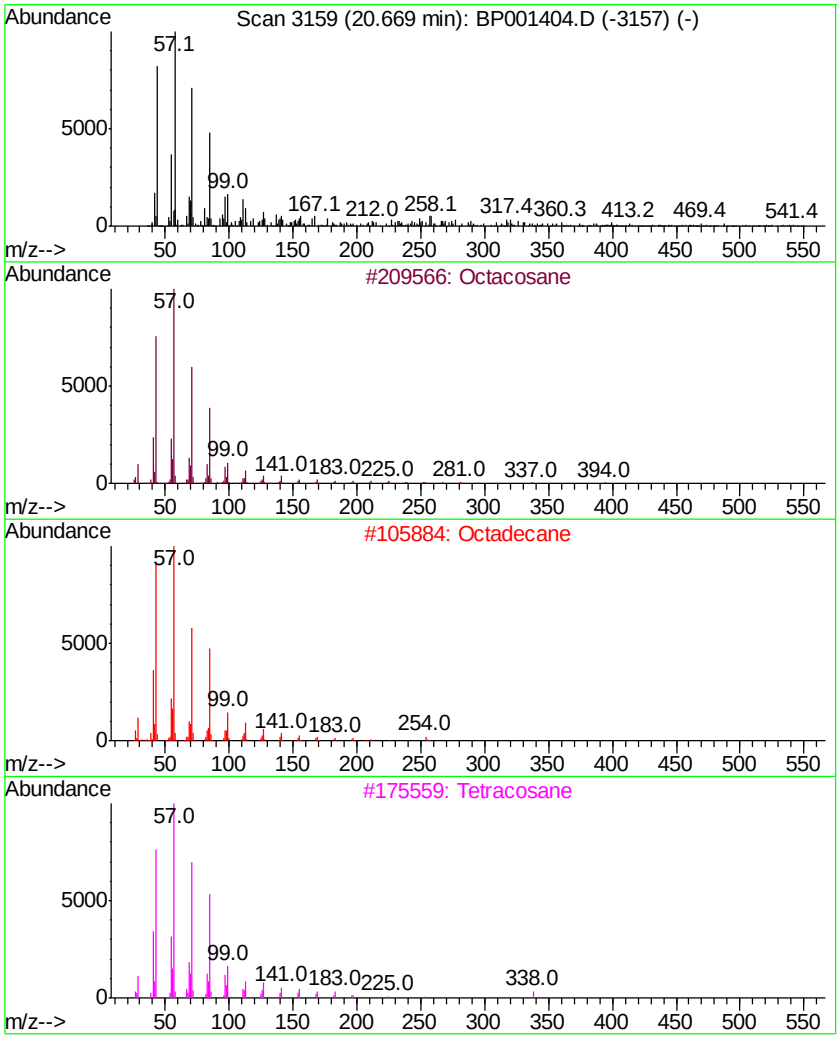
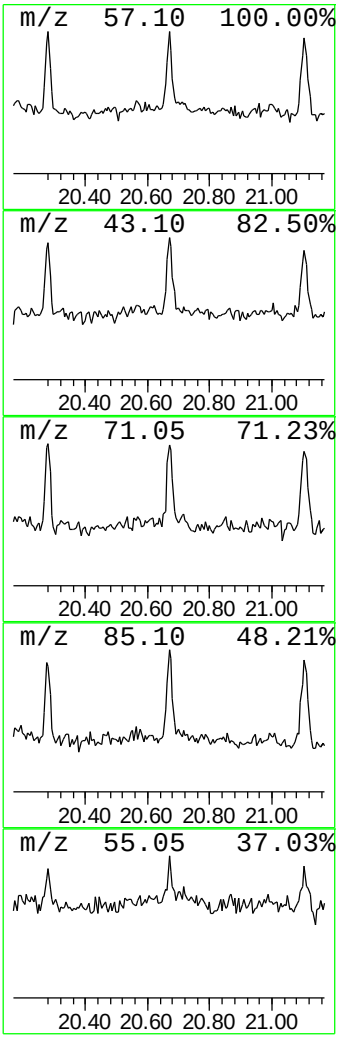
Peak Number 13 Octacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.67	3.02 ng	47071	Perylene-d12	20.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	90
2	Octadecane	254	C18H38	000593-45-3	90
3	Tetracosane	338	C24H50	000646-31-1	90
4	Eicosane	282	C20H42	000112-95-8	81
5	Heptadecane	240	C17H36	000629-78-7	81





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Operator : JU
Sample : K6413-06
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : K005-AA-WC-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP121919.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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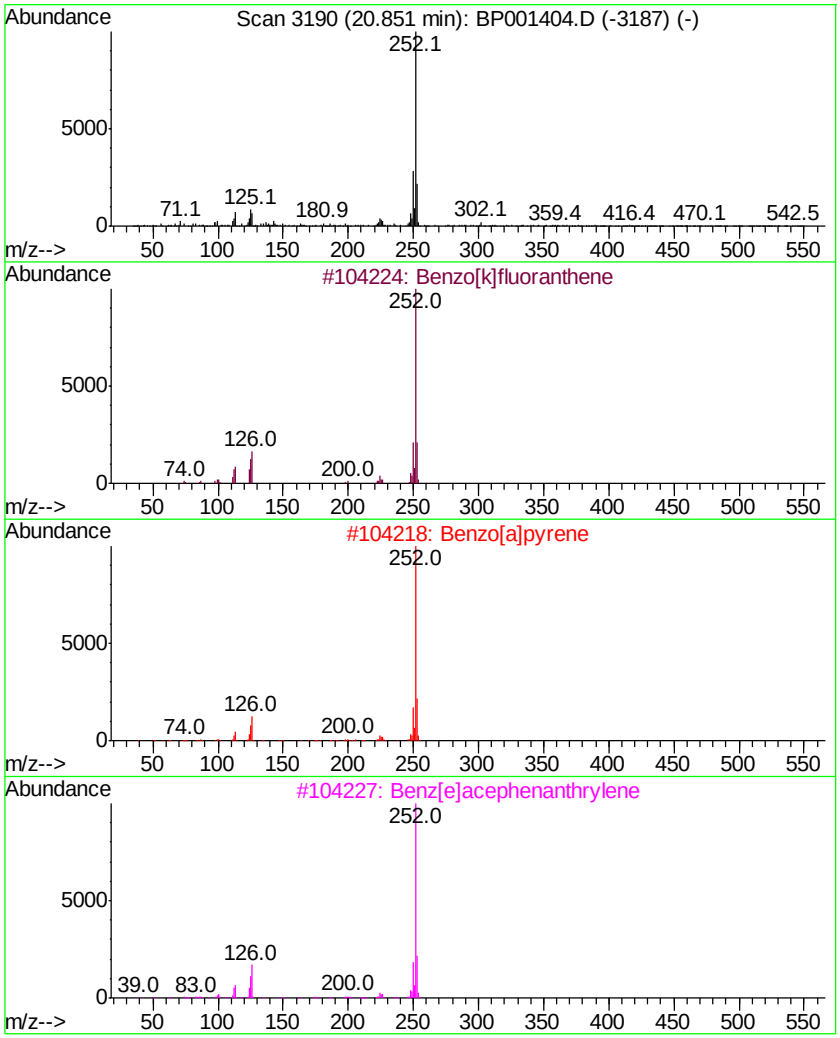
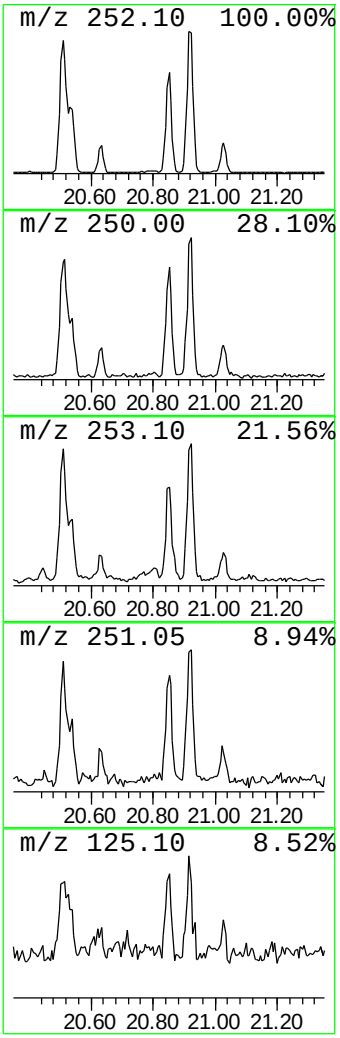
Peak Number 14 Benzo[e]pyrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.85	5.40 ng	84126	Perylene-d12	20.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
2	Benzo[a]pyrene	252	C20H12	000050-32-8	90
3	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	90
4	Perylene	252	C20H12	000198-55-0	89
5	Benzo[e]pyrene	252	C20H12	000192-97-2	81





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP122419\
Data File : BP001404.D
Acq On : 25 Dec 2019 00:04
Operator : JU
Sample : K6413-06
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : K005-AA-WC-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP121919.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...	5.60	13.9	ng	235912	1	8.29	339789	20.0
unknown7.93	7.93	76.3	ng	1295620	1	8.29	339789	20.0
Phenanthrene, 2-m...	16.33	2.9	ng	60844	4	15.58	415509	20.0
Phenanthrene, 1-m...	16.37	3.2	ng	65676	4	15.58	415509	20.0
n-Hexadecanoic acid	16.47	10.3	ng	213863	4	15.58	415509	20.0
1H-Indene, 1-phenyl-	16.52	2.0	ng	42235	4	15.58	415509	20.0
Naphthalene, 2-ph...	16.77	2.9	ng	59228	4	15.58	415509	20.0
9,10-Dimethylanth...	17.12	2.0	ng	42075	4	15.58	415509	20.0
Pyrene, 1-methyl-	18.22	2.2	ng	59758	5	19.23	544012	20.0
Cyclopentane, (4-...	19.10	6.4	ng	173870	5	19.23	544012	20.0
Tricosane	20.28	3.6	ng	56124	6	20.99	311537	20.0
Octacosane	20.67	3.0	ng	47071	6	20.99	311537	20.0
Benzo[e]pyrene	20.85	5.4	ng	84126	6	20.99	311537	20.0