LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112321\
Data File : BP008106.D
Acq On : 23 Nov 2021 20:43
Operator : CG/JU
Sample : M4782-01 2X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : CL-02-111921

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\8270E-BP111921.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP008106.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.028	3	7	24	rVB	343349	563948	15.79%	1.172%
2	4.946	328	333	341	rBV	106382	159517	4.47%	0.331%
3	5.411	406	412	427	rBV	1324746	2068887	57.94%	4.298%
4	7.005	672	683	697	rBV	1329703	2216509	62.07%	4.605%
5	7.822	814	822	831	rBV	707541	1150411	32.22%	2.390%
6	8.981	1012	1019	1029	rBV	853790	1409190	39.46%	2.928%
7	10.410	1256	1262	1270	rBV	66574	126572	3.54%	0.263%
8	10.622	1288	1298	1304	rBV	955648	1631735	45.70%	3.390%
9	10.669	1304	1306	1312	rVB	35386	50630	1.42%	0.105%
10	12.287	1575	1581	1589	rVB	53690	96371	2.70%	0.200%
11	12.504	1613	1618	1627	rVB	66207	108674	3.04%	0.226%
12	13.087	1710	1717	1724	rBV	1945730	2902757	81.29%	6.030%
13	13.298	1749	1753	1760	rVB2	35627	55154	1.54%	0.115%
14	13.645	1804	1812	1816	rBV3	42042	111578	3.12%	0.232%
15	13.787	1831	1836	1841	rBV	92379	154386	4.32%	0.321%
16	13.840	1841	1845	1850	rVB	59211	84622	2.37%	0.176%
17	14.022	1872	1876	1880	rVV	41808	57384	1.61%	0.119%
18	14.187	1899	1904	1913	rBV2	106753	185337	5.19%	0.385%
19	14.298	1919	1923	1932	rVB5	37119	69361	1.94%	0.144%
20	14.375	1932	1936	1940	rVB	27111	36126	1.01%	0.075%
21	14.469	1943	1952	1957	rVV	1444510	2198180	61.56%	4.567%
22	14.528	1958	1962	1971	rVB	121890	199029	5.57%	0.413%
23	14.687	1983	1989	1996	rVB4	25868	66054	1.85%	0.137%
24	14.869	2015	2020	2026	rVB	107763	169605	4.75%	0.352%
25	14.945	2029	2033	2038	rVB2	60701	83544	2.34%	0.174%
26	15.092	2051	2058	2062	rBV4	56410	116953	3.28%	0.243%
27	15.139	2062	2066	2070	rVB	42610	63255	1.77%	0.131%
28	15.257	2079	2086	2089	rBV5	59811	131333	3.68%	0.273%
29	15.292	2089	2092	2096	rVV	63123	91398	2.56%	0.190%
30	15.334	2096	2099	2105	rVB2	25265	43218	1.21%	0.090%
31	15.516	2125	2130	2135	rBV2	217262	326562	9.15%	0.678%
32	15.645	2148	2152	2155	rBV3	30918	40956	1.15%	0.085%
33	15.681	2155	2158	2162	rVB4	34425	47125	1.32%	0.098%
34	15.734	2163	2167	2175	rBV7	37465	73264	2.05%	0.152%
35	15.816	2177	2181	2190	rVB2	56100	110886	3.11%	0.230%
36	15.963	2199	2206	2214	rBV	1656704	2468528	69.13%	5.128%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112321\
Data File : BP008106.D
Acq On : 23 Nov 2021 20:43
Operator : CG/JU
Sample : M4782-01 2X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CL-02-111921

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\8270E-BP111921.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	16.198	2242	2246	2257	rVB4	76167	159543	4.47%	0.331%
38	16.339	2265	2270	2273	rBV2	32363	52746	1.48%	0.110%
39	16.528	2295	2302	2306	rBV2	79275	160772	4.50%	0.334%
40	16.575	2307	2310	2316	rVB3	73126	100936	2.83%	0.210%
41	16.681	2325	2328	2332	rVB2	51737	68742	1.93%	0.143%
42	16.763	2339	2342	2345	rBV4	29751	46488	1.30%	0.097%
43	16.881	2359	2362	2368	rVB3	37170	62117	1.74%	0.129%
44	16.957	2372	2375	2379	rBV5	27216	47770	1.34%	0.099%
45	17.039	2385	2389	2399	rVB2	122276	217039	6.08%	0.451%
46	17.228	2415	2421	2424	rBV	1729181	2486058	69.62%	5.165%
47	17.269	2424	2428	2434	rVV	1457725	2089697	58.52%	4.341%
48	17.357	2438	2443	2449	rVB	476024	684637	19.17%	1.422%
49	17.422	2449	2454	2459	rBV2	65779	133643	3.74%	0.278%
50	17.633	2487	2490	2500	rBV3	69085	129445	3.62%	0.269%
51	17.810	2516	2520	2525	rVB	76878	101403	2.84%	0.211%
52	18.092	2564	2568	2571	rBV	309511	445058	12.46%	0.925%
53	18.145	2573	2577	2583	rVV	444765	646313	18.10%	1.343%
54	18.222	2586	2590	2595	rVV	171945	249477	6.99%	0.518%
55	18.281	2595	2600	2604	rVV2	502631	886953	24.84%	1.843%
56	18.322	2604	2607	2612	rVB	268063	343021	9.61%	0.713%
57	18.580	2647	2651	2655	rBV	179186	218324	6.11%	0.454%
58	18.822	2689	2692	2695	rVB	75452	80854	2.26%	0.168%
59	18.869	2697	2700	2703	rVV	94908	102047	2.86%	0.212%
60	18.986	2716	2720	2724	rBV	280044	436169	12.21%	0.906%
61	19.122	2740	2743	2750	rBV4	116681	252457	7.07%	0.524%
62	19.245	2758	2764	2770	rBV	2149663	2887019	80.85%	5.998%
63	19.592	2814	2823	2832	rBV	1851789	3314092	92.81%	6.885%
64	19.792	2852	2857	2861	rVB	2701630	3277592	91.79%	6.809%
65	19.910	2874	2877	2884	rVB2	139314	287893	8.06%	0.598%
66	20.080	2903	2906	2913	rVB	446711	584099	16.36%	1.213%
67	20.175	2919	2922	2926	rVB2	328452	390819	10.94%	0.812%
68	20.233	2930	2932	2936	rVB	248606	274105	7.68%	0.569%
69	20.375	2953	2956	2959	rBV	159441	179167	5.02%	0.372%
70	20.416	2960	2963	2966	rBV	187629	218721	6.13%	0.454%
71	21.322	3111	3117	3121	rBV2	1798975	3570921	100.00%	7.419%
72	21.363	3121	3124	3134	rVB	892995	1251165	35.04%	2.599%
73	22.998	3399	3402	3407	rBV	375837	614289	17.20%	1.276%
74	23.733	3522	3527	3532	rVB2	861417	1614707	45.22%	3.355%

Sum of corrected areas: 48135337

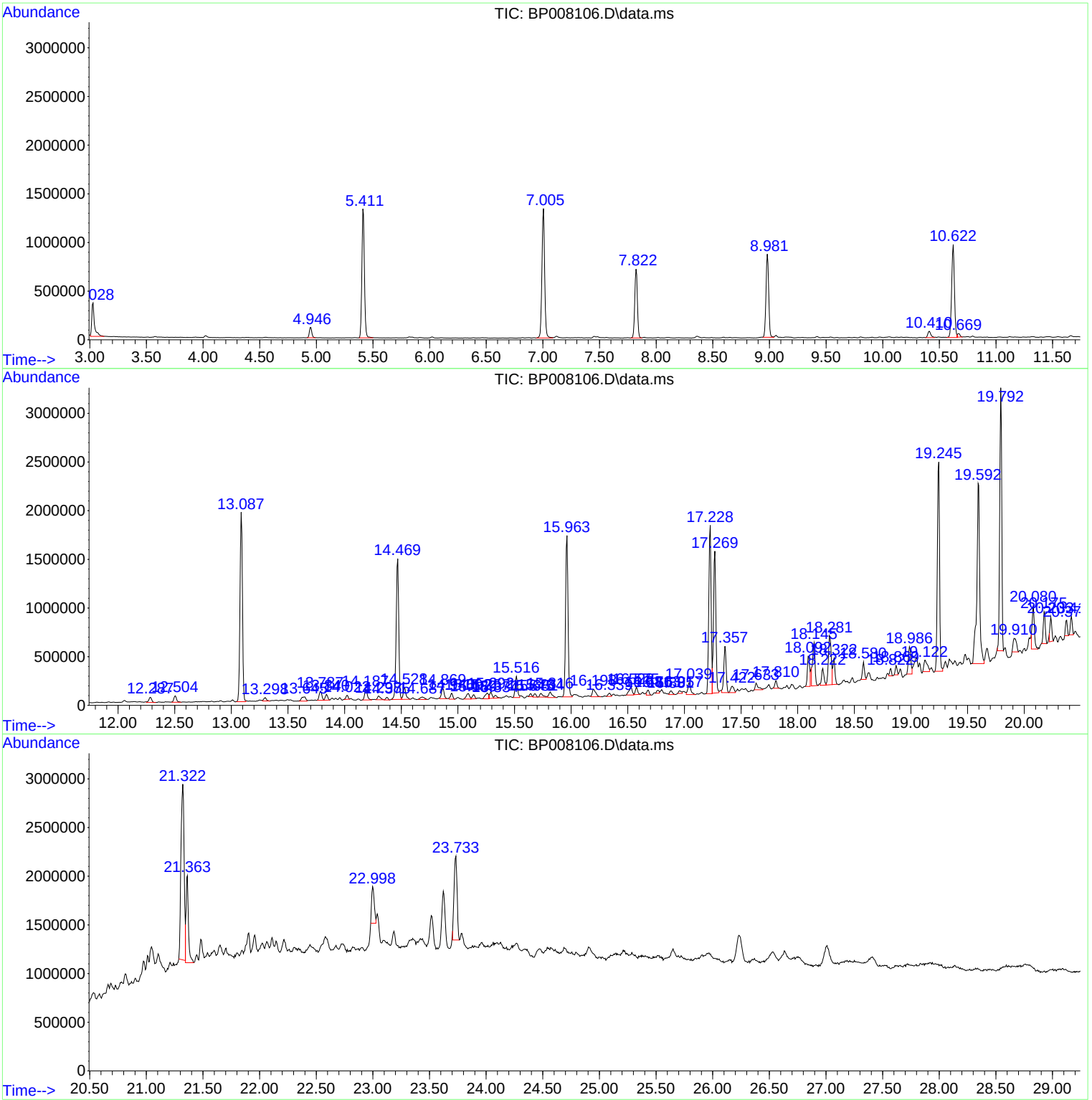
Instrument :
BNA_P
ClientSampleId :
CL-02-111921

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112321\
Data File : BP008106.D
Acq On : 23 Nov 2021 20:43
Operator : CG/JU
Sample : M4782-01 2X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CL-02-111921

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112321\
Data File : BP008106.D
Acq On : 23 Nov 2021 20:43
Operator : CG/JU
Sample : M4782-01 2X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CL-02-111921

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

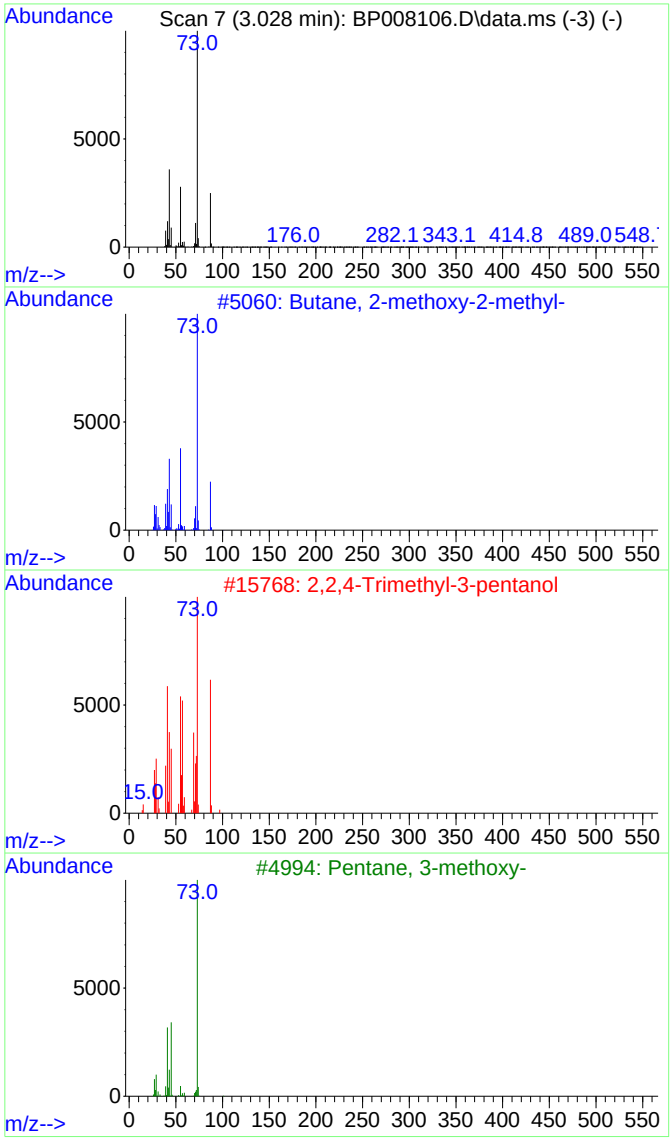
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

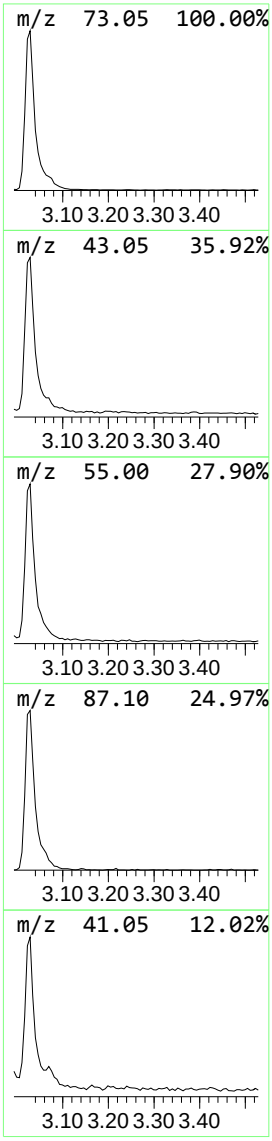
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.028	9.80 ng	563948	1,4-Dichlorobenzene-d4	7.822

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2	2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	40
3	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
4	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
5	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112321\
Data File : BP008106.D
Acq On : 23 Nov 2021 20:43
Operator : CG/JU
Sample : M4782-01 2X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CL-02-111921

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

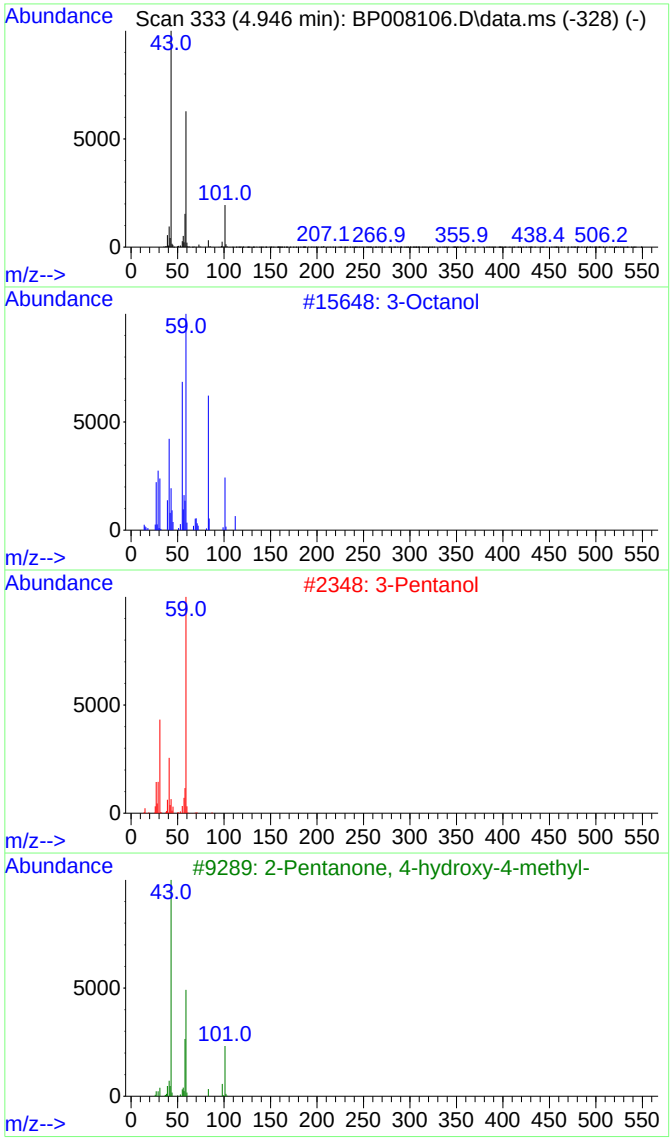
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

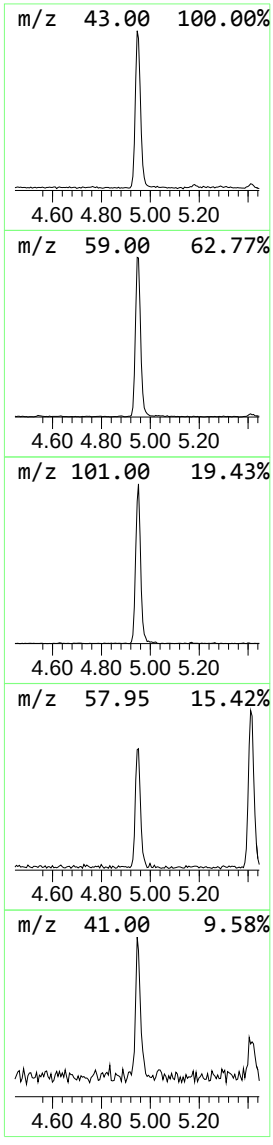
****

Peak Number 2 3-Octanol Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.946	2.77 ng	159517	1,4-Dichlorobenzene-d4	7.822

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Octanol	130	C8H18O	000589-98-0	38
2	3-Pentanol	88	C5H12O	000584-02-1	32
3	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	32
4	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	25
5	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112321\
Data File : BP008106.D
Acq On : 23 Nov 2021 20:43
Operator : CG/JU
Sample : M4782-01 2X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CL-02-111921

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

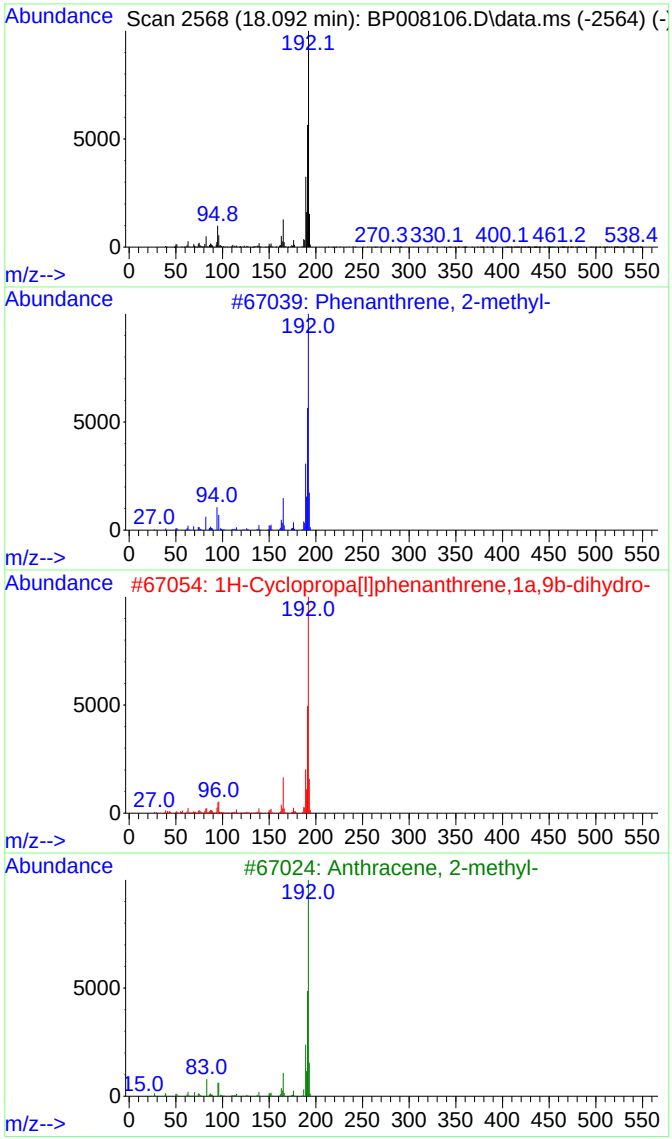
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

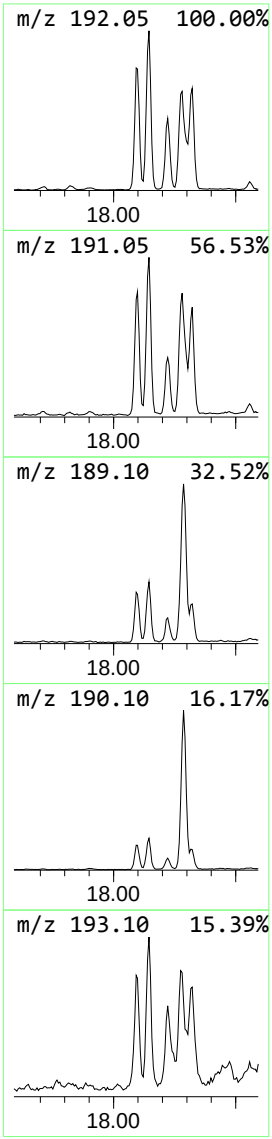
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.092	3.58 ng	445058	Phenanthrene-d10	17.228

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	96
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	95
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112321\
Data File : BP008106.D
Acq On : 23 Nov 2021 20:43
Operator : CG/JU
Sample : M4782-01 2X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CL-02-111921

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

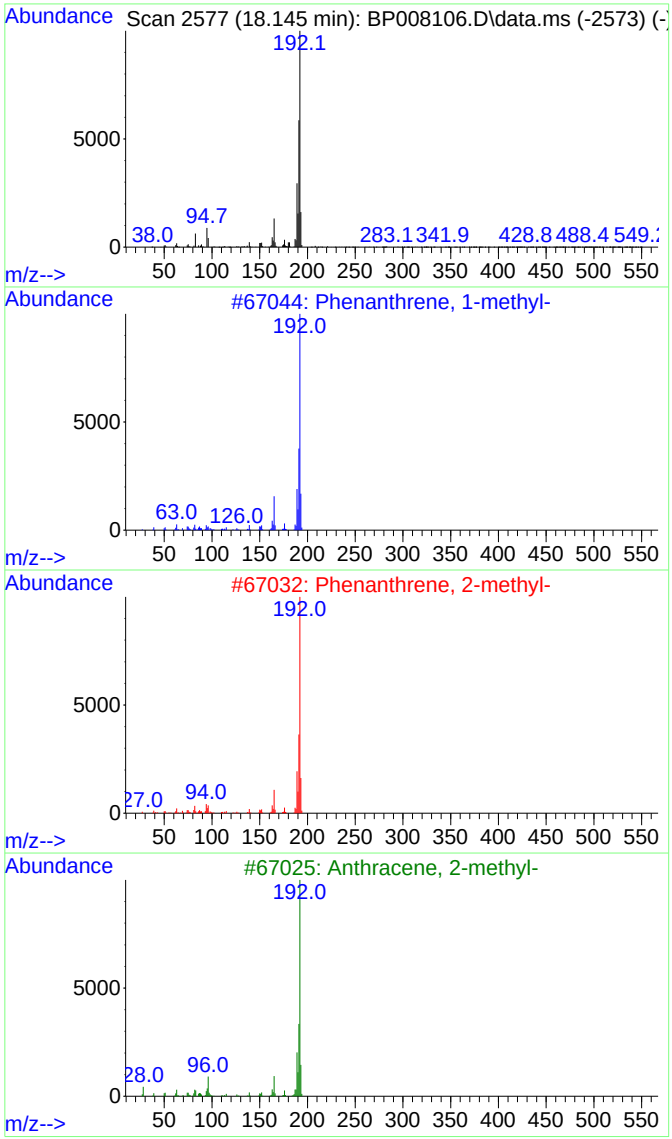
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

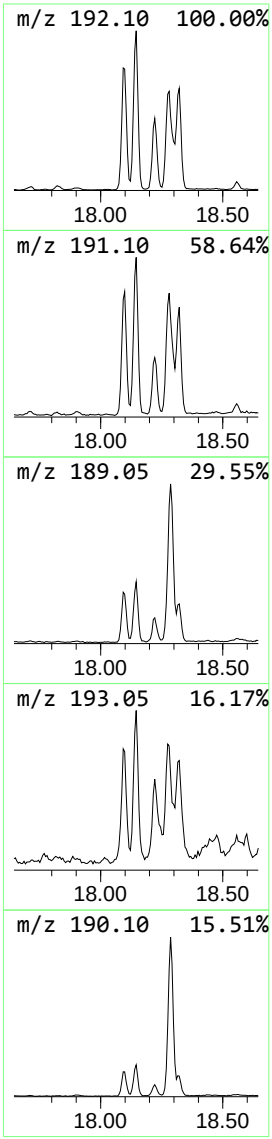
****

Peak Number 5 Phenanthrene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.145	5.20 ng	646313	Phenanthrene-d10	17.228

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
4	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	93
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112321\
Data File : BP008106.D
Acq On : 23 Nov 2021 20:43
Operator : CG/JU
Sample : M4782-01 2X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CL-02-111921

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

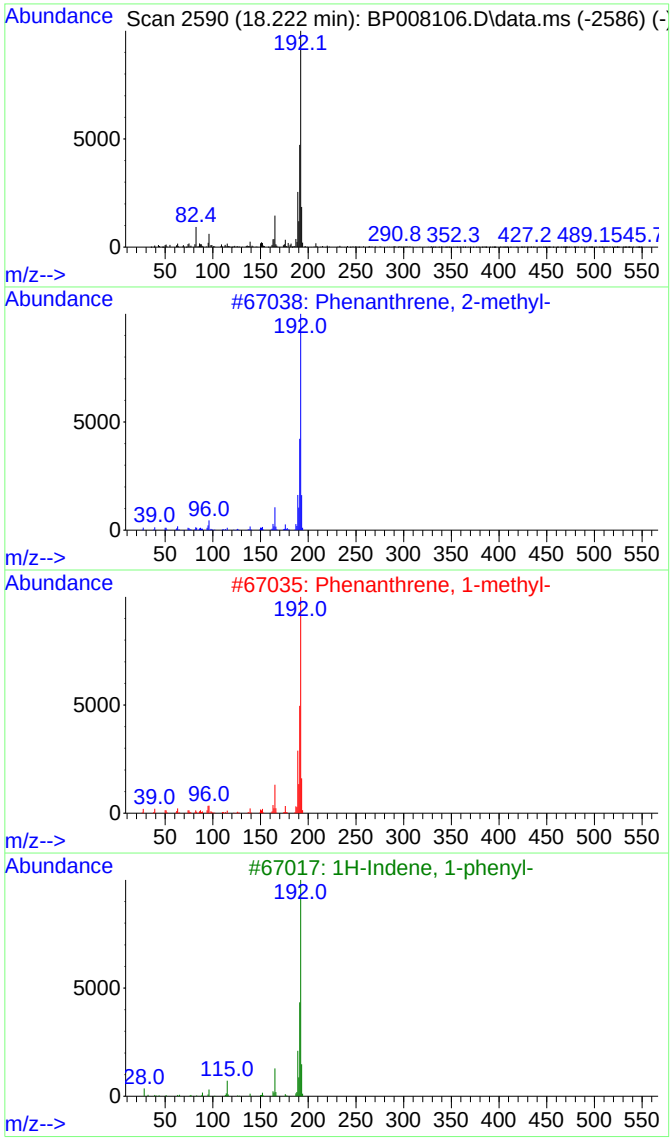
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

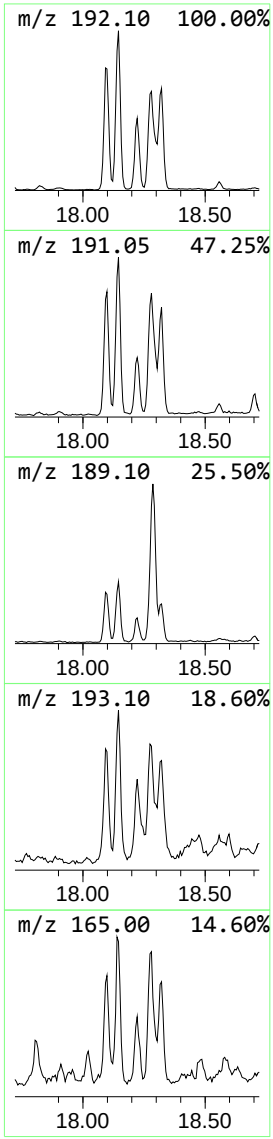
****

Peak Number 6 1H-Indene, 1-phenyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.222	2.01 ng	249477	Phenanthrene-d10	17.228

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94
3	1H-Indene, 1-phenyl-	192	C15H12	001961-96-2	93
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
5	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112321\
Data File : BP008106.D
Acq On : 23 Nov 2021 20:43
Operator : CG/JU
Sample : M4782-01 2X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CL-02-111921

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

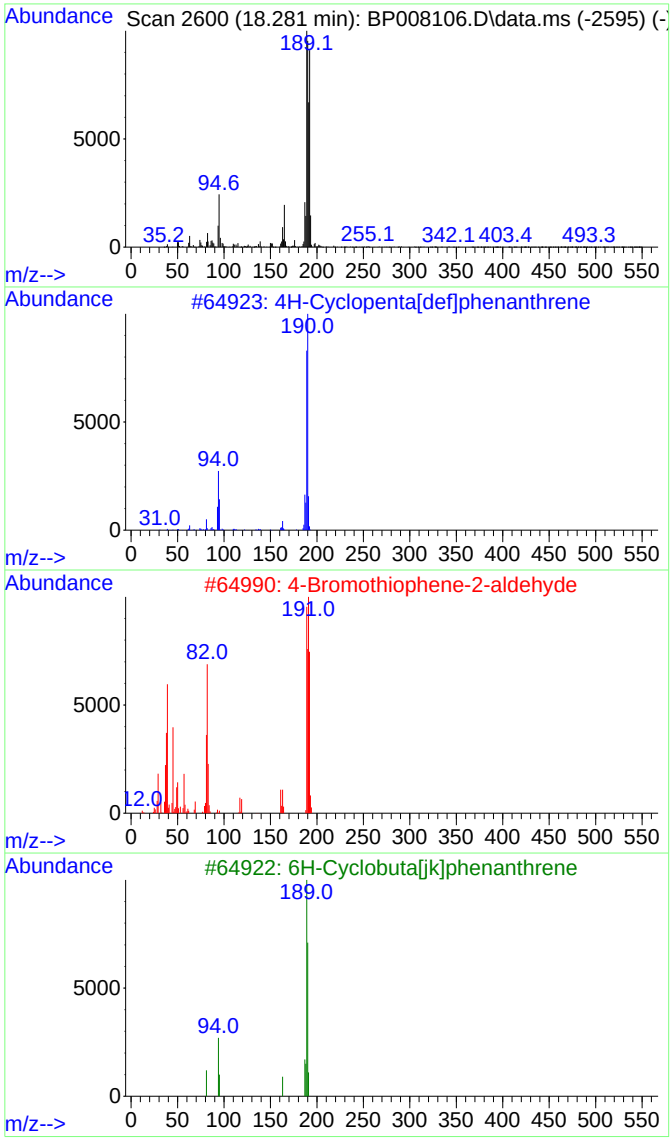
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

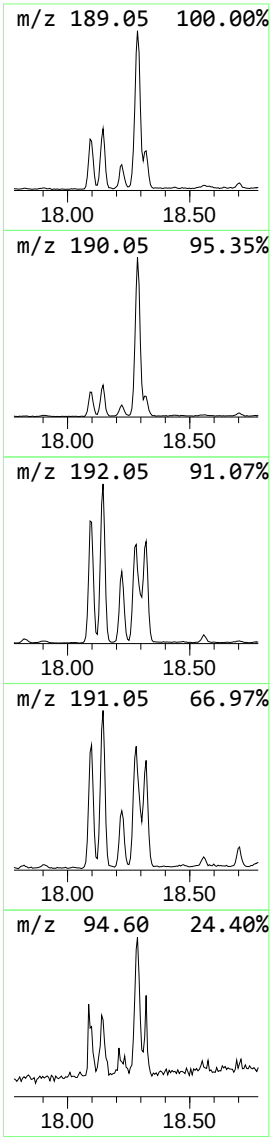
****

Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.281	7.14 ng	886953	Phenanthrene-d10	17.228

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	60
2	4-Bromothiophene-2-aldehyde	190	C5H3BrOS	018791-75-8	58
3	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	49
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	46
5	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	46





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112321\
Data File : BP008106.D
Acq On : 23 Nov 2021 20:43
Operator : CG/JU
Sample : M4782-01 2X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CL-02-111921

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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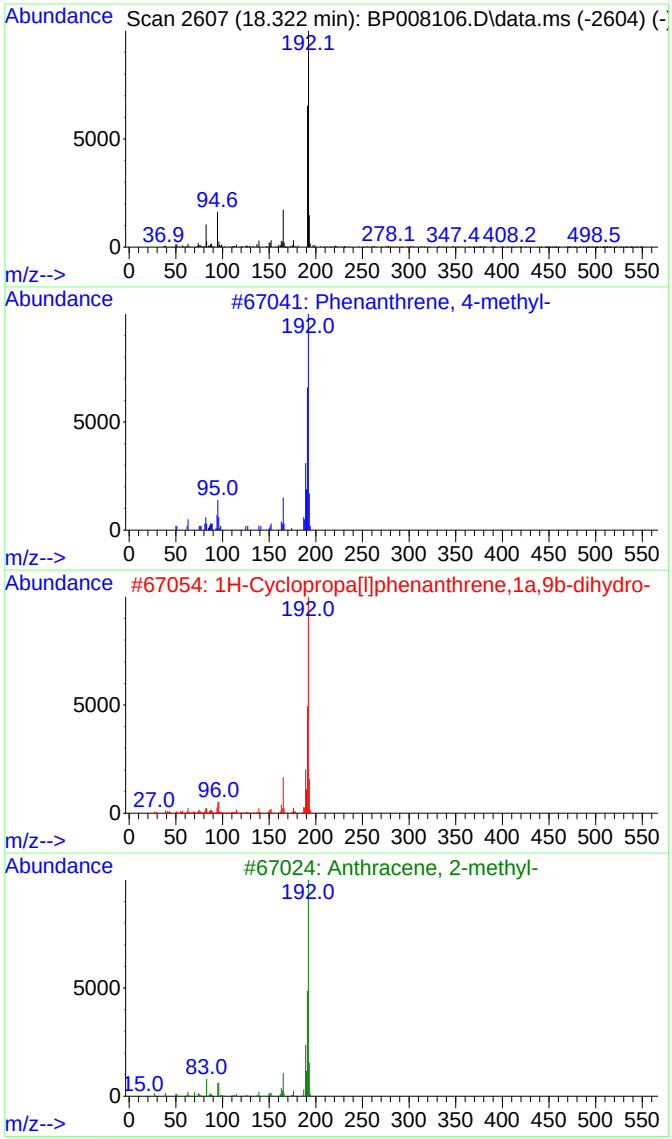
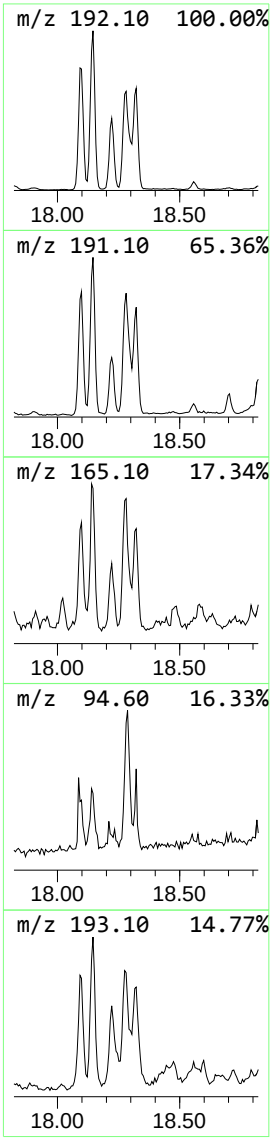
Peak Number 8 Phenanthrene, 4-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.322	2.76 ng	343021	Phenanthrene-d10	17.228

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	87
2	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	81
3	Anthracene, 2-methyl-	192	C15H12	000613-12-7	81
4	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	81
5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	81





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Data File : BP008106.D
Acq On : 23 Nov 2021 20:43
Operator : CG/JU
Sample : M4782-01 2X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CL-02-111921

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

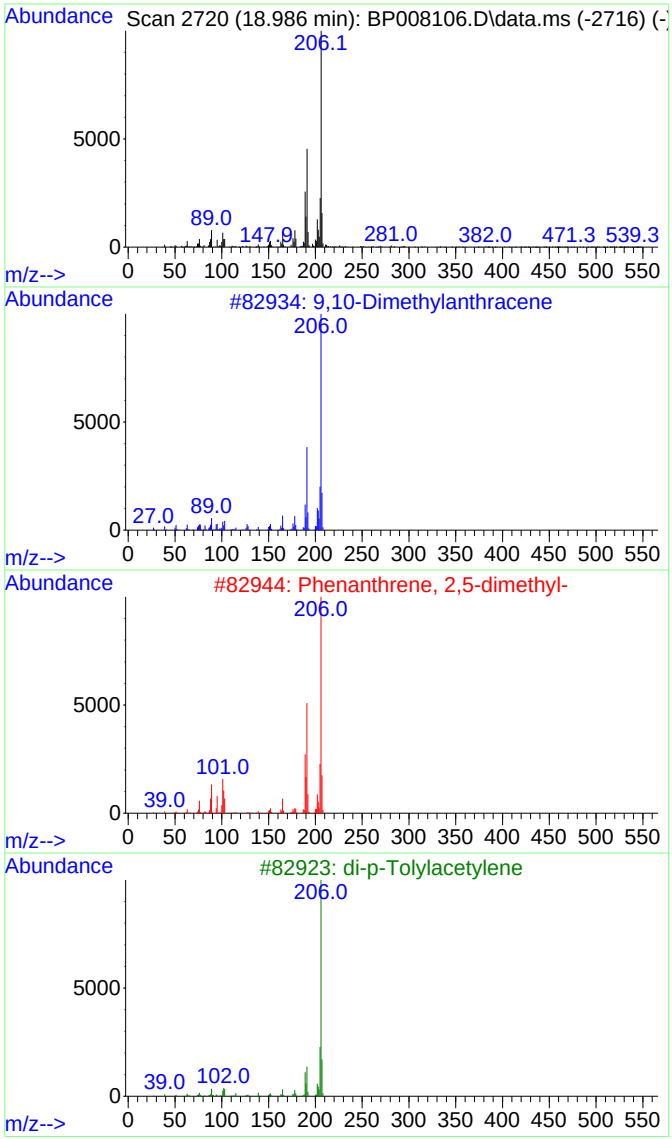
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

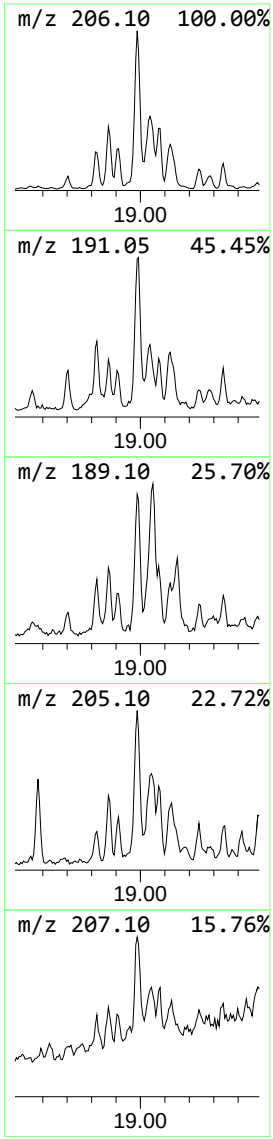
****

Peak Number 9 9,10-Dimethylanthracene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.986	3.51 ng	436169	Phenanthrene-d10	17.228

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Dimethylanthracene	206	C16H14	000781-43-1	98
2	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	95
3	di-p-Tolylacetylene	206	C16H14	002789-88-0	93
4	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	91
5	Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	89





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Data File : BP008106.D
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Operator : CG/JU
Sample : M4782-01 2X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CL-02-111921

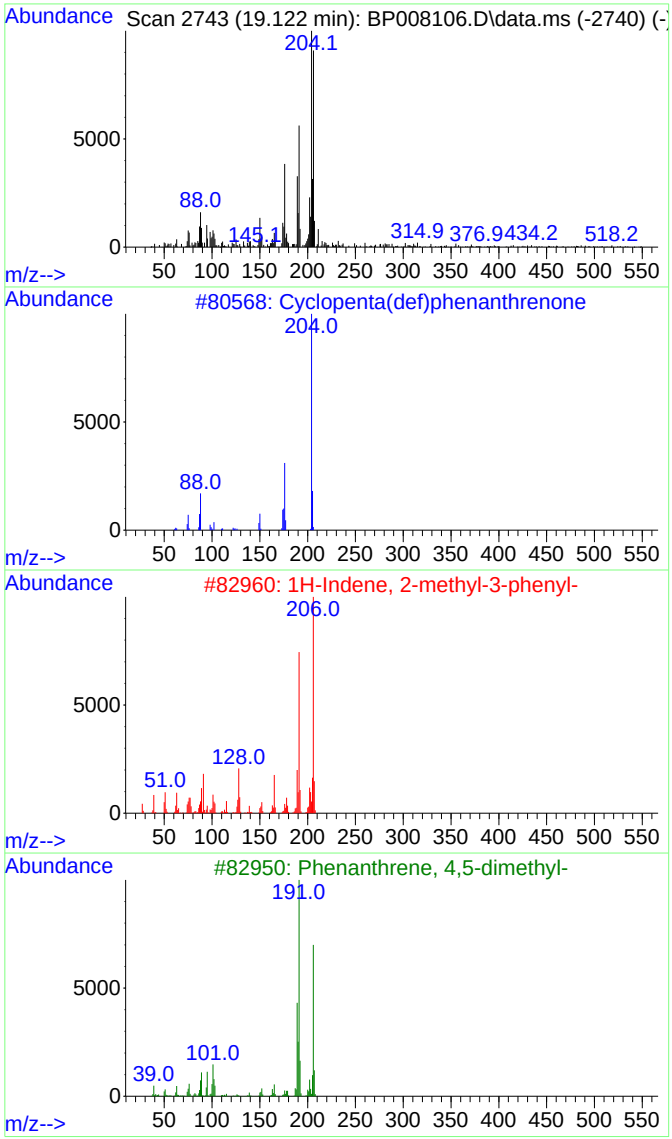
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

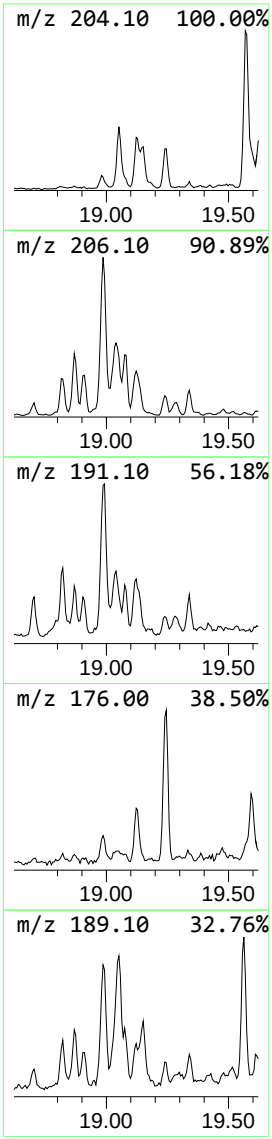
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Cyclopenta(def)phenanthrenone Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.122	2.03 ng	252457	Phenanthrene-d10	17.228

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	86
2	1H-Indene, 2-methyl-3-phenyl-	206	C16H14	035099-60-6	46
3	Phenanthrene, 4,5-dimethyl-	206	C16H14	003674-69-9	46
4	9,10-Dimethylanthracene	206	C16H14	000781-43-1	42
5	1,3-Diphenyl-3-methylcyclopropene	206	C16H14	086544-79-8	41





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Data File : BP008106.D
Acq On : 23 Nov 2021 20:43
Operator : CG/JU
Sample : M4782-01 2X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CL-02-111921

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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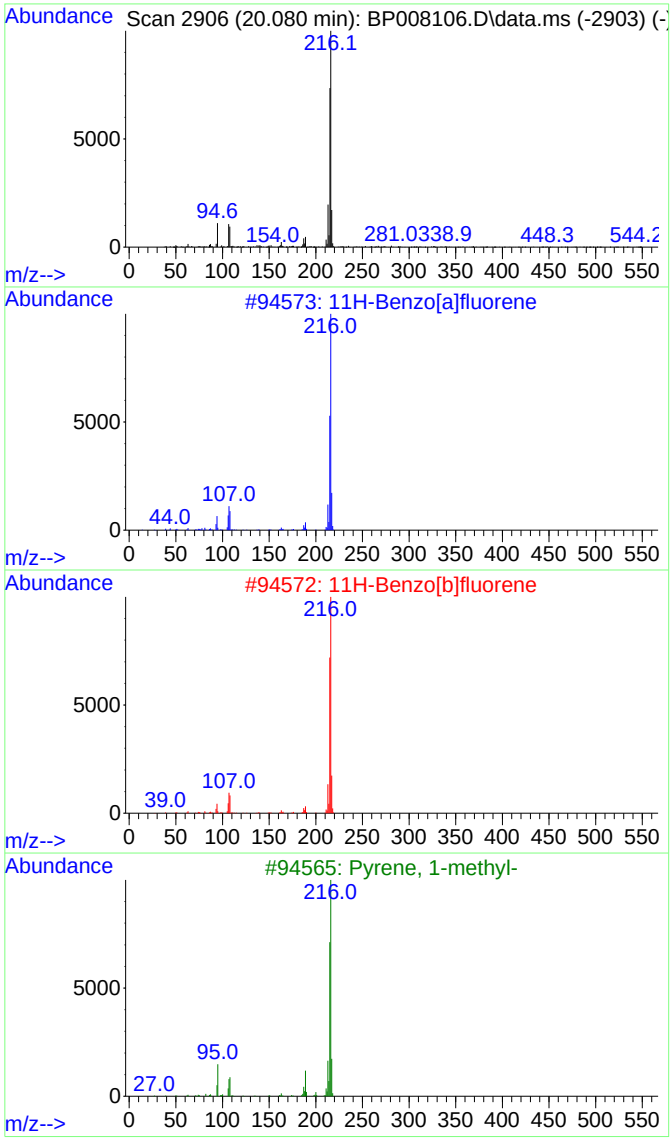
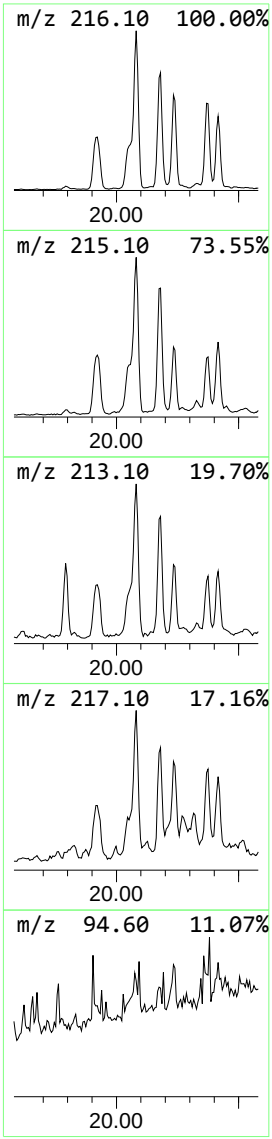
Peak Number 11 11H-Benzo[a]fluorene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.080	3.27 ng	584099	Chrysene-d12	21.327

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	95
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
3	Pyrene, 1-methyl-	216	C17H12	002381-21-7	90
4	Pyrene, 2-methyl-	216	C17H12	003442-78-2	87
5	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112321\
Data File : BP008106.D
Acq On : 23 Nov 2021 20:43
Operator : CG/JU
Sample : M4782-01 2X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CL-02-111921

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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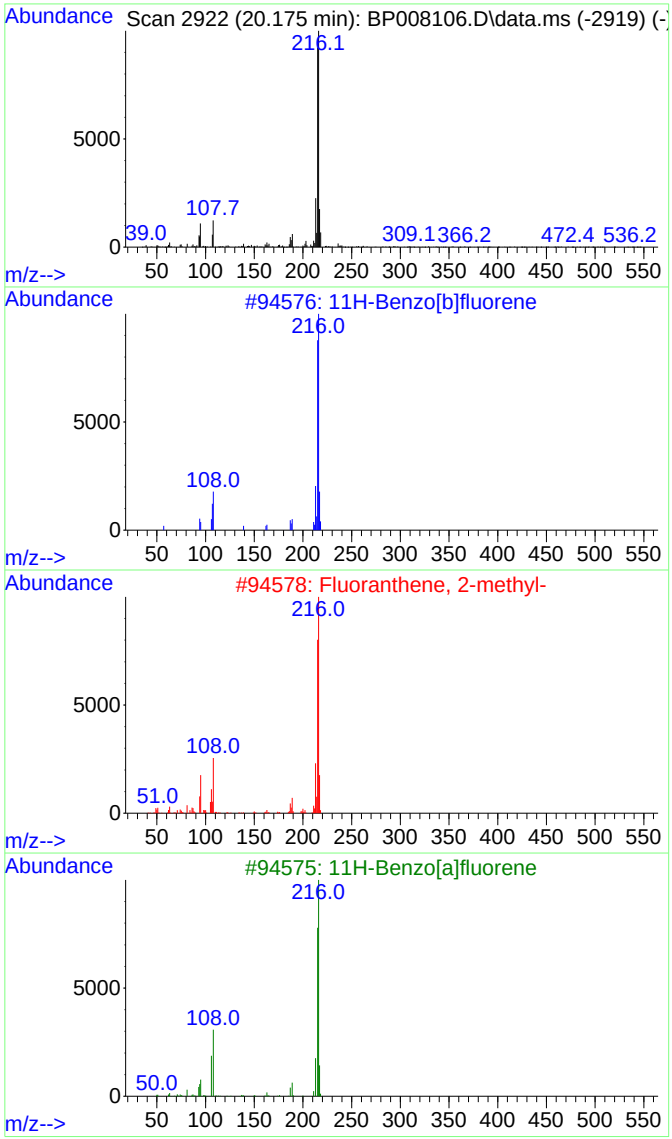
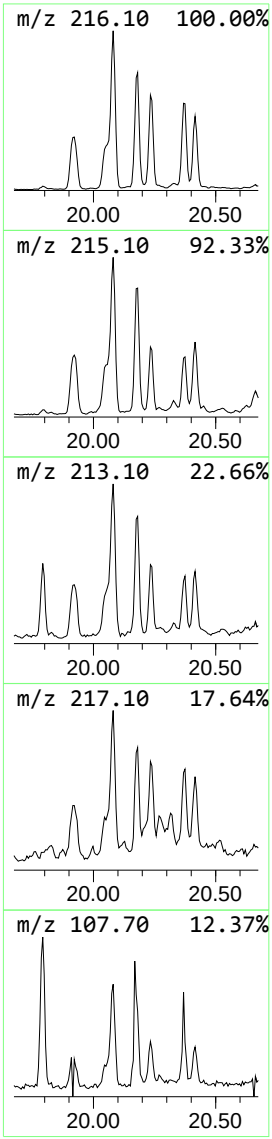
Peak Number 12 11H-Benzo[b]fluorene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.174	2.19 ng	390819	Chrysene-d12	21.327

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	95
2	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	93
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	93
4	Pyrene, 1-methyl-	216	C17H12	002381-21-7	89
5	6H-Thiazolo[5,4-e]indole, 2,7,8-...	216	C12H12N2S	1000338-32-0	72





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112321\
Data File : BP008106.D
Acq On : 23 Nov 2021 20:43
Operator : CG/JU
Sample : M4782-01 2X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CL-02-111921

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-metho...	3.028	9.8	ng	563948	1	7.822	1150410	20.0
3-Octanol	4.946	2.8	ng	159517	1	7.822	1150410	20.0
Phenanthrene, 2...	18.092	3.6	ng	445058	4	17.228	2486060	20.0
Phenanthrene, 1...	18.145	5.2	ng	646313	4	17.228	2486060	20.0
1H-Indene, 1-ph...	18.222	2.0	ng	249477	4	17.228	2486060	20.0
4H-Cyclopenta[d...	18.281	7.1	ng	886953	4	17.228	2486060	20.0
Phenanthrene, 4...	18.322	2.8	ng	343021	4	17.228	2486060	20.0
9,10-Dimethylan...	18.986	3.5	ng	436169	4	17.228	2486060	20.0
Cyclopenta(def)...	19.122	2.0	ng	252457	4	17.228	2486060	20.0
11H-Benzo[a]flu...	20.080	3.3	ng	584099	5	21.327	3570920	20.0
11H-Benzo[b]flu...	20.174	2.2	ng	390819	5	21.327	3570920	20.0