LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062824\
Data File : BM046352.D
Acq On : 28 Jun 2024 14:44
Operator : MA/JU
Sample : P3047-01 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EO-01-062424

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_M\Methods\8270-BM062724.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM046352.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.610	289	293	301	rBV2	10107	23347	1.37%	0.108%
2	5.045	361	367	385	rBV	193182	409392	24.00%	1.897%
3	6.639	630	638	657	rBV	145852	410818	24.08%	1.904%
4	7.375	755	763	782	rBV	553984	1040613	61.00%	4.822%
5	7.922	849	856	863	rBV9	8559	22276	1.31%	0.103%
6	8.563	959	965	979	rBV	81320	236666	13.87%	1.097%
7	10.133	1224	1232	1257	rBV	607540	1460052	85.59%	6.766%
8	11.827	1513	1520	1537	rBV2	43089	121085	7.10%	0.561%
9	12.039	1548	1556	1569	rBV	51481	116079	6.80%	0.538%
10	12.645	1653	1659	1678	rBV	265250	525322	30.80%	2.434%
11	13.204	1745	1754	1760	rBV	37387	103544	6.07%	0.480%
12	13.339	1771	1777	1782	rBV	74100	136813	8.02%	0.634%
13	13.392	1782	1786	1796	rVB	47326	89623	5.25%	0.415%
14	13.574	1813	1817	1821	rBV	32094	51868	3.04%	0.240%
15	13.610	1821	1823	1828	rVB2	16477	18799	1.10%	0.087%
16	13.745	1839	1846	1855	rBV	91224	163070	9.56%	0.756%
17	14.021	1883	1893	1901	rBV	1034420	1617555	94.82%	7.496%
18	14.086	1901	1904	1910	rVB2	34241	54682	3.21%	0.253%
19	14.257	1927	1933	1936	rBV6	14708	34411	2.02%	0.159%
20	14.439	1955	1964	1972	rVB3	29165	63691	3.73%	0.295%
21	14.515	1972	1977	1983	rBV	34734	55452	3.25%	0.257%
22	14.651	1994	2000	2006	rBV3	44595	87586	5.13%	0.406%
23	14.710	2006	2010	2016	rVB3	20963	34238	2.01%	0.159%
24	14.827	2023	2030	2033	rBV4	25007	51888	3.04%	0.240%
25	14.857	2033	2035	2039	rVB2	15775	17982	1.05%	0.083%
26	14.904	2039	2043	2048	rVB3	25322	32617	1.91%	0.151%
27	14.992	2052	2058	2060	rBV4	15795	30089	1.76%	0.139%
28	15.027	2060	2064	2068	rVB2	21931	32761	1.92%	0.152%
29	15.086	2068	2074	2080	rBV3	81889	166324	9.75%	0.771%
30	15.245	2097	2101	2105	rVB5	13427	21703	1.27%	0.101%
31	15.298	2105	2110	2116	rVV6	16535	34949	2.05%	0.162%
32	15.386	2119	2125	2134	rVB6	19648	49522	2.90%	0.229%
33	15.533	2140	2150	2158	rBV3	202512	395747	23.20%	1.834%
34	15.768	2185	2190	2200	rBV9	28017	63195	3.70%	0.293%
35	16.086	2236	2244	2249	rBV	68137	151794	8.90%	0.703%
36	16.133	2249	2252	2258	rVB	51155	73466	4.31%	0.340%

LSC Area Percent Report

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Acq On : 28 Jun 2024 14:44
Operator : MA/JU
Sample : P3047-01 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EO-01-062424

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_M\Methods\8270-BM062724.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	16.233	2265	2269	2275	rVB3	53224	77694	4.55%	0.360%
38	16.321	2281	2284	2286	rVV4	13409	18792	1.10%	0.087%
39	16.356	2287	2290	2295	rVB5	21827	31630	1.85%	0.147%
40	16.439	2300	2304	2311	rVB6	24369	47273	2.77%	0.219%
41	16.515	2312	2317	2321	rBV6	17978	34819	2.04%	0.161%
42	16.592	2326	2330	2341	rVB2	54819	119977	7.03%	0.556%
43	16.774	2354	2361	2365	rBV	1132603	1696757	99.47%	7.863%
44	16.815	2365	2368	2377	rVV	441355	654216	38.35%	3.032%
45	16.909	2380	2384	2390	rVB	94995	146314	8.58%	0.678%
46	16.974	2390	2395	2400	rBV2	34704	71531	4.19%	0.331%
47	17.133	2419	2422	2425	rVB4	14600	18254	1.07%	0.085%
48	17.198	2428	2433	2435	rBV4	17163	32782	1.92%	0.152%
49	17.298	2446	2450	2453	rVB3	13523	20175	1.18%	0.093%
50	17.356	2456	2460	2465	rVB2	50221	72964	4.28%	0.338%
51	17.498	2480	2484	2492	rVB	39201	79710	4.67%	0.369%
52	17.574	2492	2497	2499	rBV5	11825	21183	1.24%	0.098%
53	17.645	2505	2509	2514	rBV	186664	279817	16.40%	1.297%
54	17.698	2514	2518	2527	rVV2	315095	509897	29.89%	2.363%
55	17.780	2527	2532	2536	rVV	94388	158338	9.28%	0.734%
56	17.839	2536	2542	2546	rVV2	254327	483675	28.35%	2.241%
57	17.880	2546	2549	2560	rVB	191747	287364	16.85%	1.332%
58	17.974	2561	2565	2570	rBV5	30162	45380	2.66%	0.210%
59	18.045	2574	2577	2581	rVB3	27055	34187	2.00%	0.158%
60	18.115	2586	2589	2591	rBV3	12982	18642	1.09%	0.086%
61	18.156	2592	2596	2600	rBV	56583	79861	4.68%	0.370%
62	18.374	2629	2633	2635	rBV	30018	43243	2.53%	0.200%
63	18.398	2635	2637	2643	rVV	44945	64493	3.78%	0.299%
64	18.456	2643	2647	2650	rVV	76134	98327	5.76%	0.456%
65	18.492	2650	2653	2658	rVB2	51124	72662	4.26%	0.337%
66	18.574	2662	2667	2671	rBV	200206	305684	17.92%	1.417%
67	18.645	2671	2679	2681	rVV5	158651	359527	21.08%	1.666%
68	18.668	2681	2683	2687	rVB	76341	87361	5.12%	0.405%
69	18.727	2687	2693	2699	rBV5	61170	156508	9.17%	0.725%
70	18.839	2707	2712	2721	rBV	434997	679432	39.83%	3.148%
71	18.945	2728	2730	2734	rVB3	37101	41938	2.46%	0.194%
72	18.992	2735	2738	2742	rBV3	86706	115243	6.76%	0.534%
73	19.086	2751	2754	2758	rBV3	28509	37895	2.22%	0.176%
74	19.203	2767	2774	2784	rVV	585964	912706	53.50%	4.229%
75	19.292	2785	2789	2795	rVB2	67441	101312	5.94%	0.469%
76	19.421	2807	2811	2822	rVB	487115	699894	41.03%	3.243%

LSC Area Percent Report

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Operator : MA/JU
Sample : P3047-01 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
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EO-01-062424

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
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If leading or trailing edge < 100 prefer < Baseline drop else tangent >
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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM062724.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	19.544	2826	2832	2842	rBV6	88961	249340	14.62%	1.155%
78	19.627	2842	2846	2849	rBV5	37658	54735	3.21%	0.254%
79	19.686	2850	2856	2857	rVV3	76224	135836	7.96%	0.629%
80	19.709	2857	2860	2864	rVB	159151	210015	12.31%	0.973%
81	19.809	2874	2877	2882	rVB	120116	147001	8.62%	0.681%
82	19.868	2883	2887	2892	rBV	151217	190252	11.15%	0.882%
83	20.009	2907	2911	2915	rBV	116811	149287	8.75%	0.692%
84	20.056	2915	2919	2923	rVV	118411	159532	9.35%	0.739%
85	20.662	3020	3022	3025	rVB	50571	50177	2.94%	0.233%
86	20.991	3071	3078	3082	rBV2	1070571	1705852	100.00%	7.905%
87	21.027	3082	3084	3096	rVB	232895	366891	21.51%	1.700%
88	21.615	3182	3184	3190	rVB2	100905	124942	7.32%	0.579%
89	23.662	3525	3532	3541	rVB	588123	1219341	71.48%	5.650%

Sum of corrected areas: 21579677

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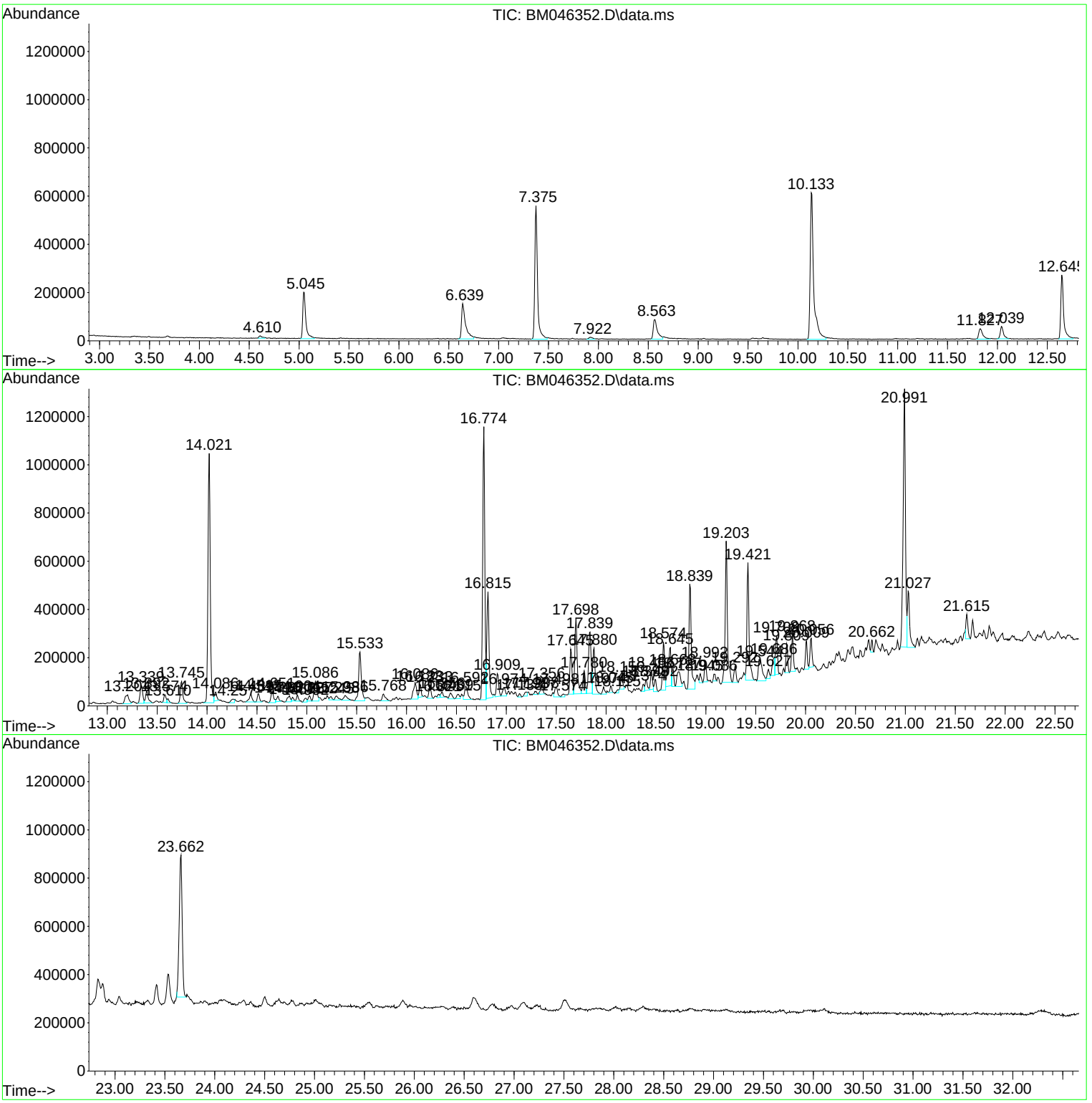
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EO-01-062424

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM062724.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_M\Data\BM062824\
Data File : BM046352.D
Acq On : 28 Jun 2024 14:44
Operator : MA/JU
Sample : P3047-01 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM062724.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

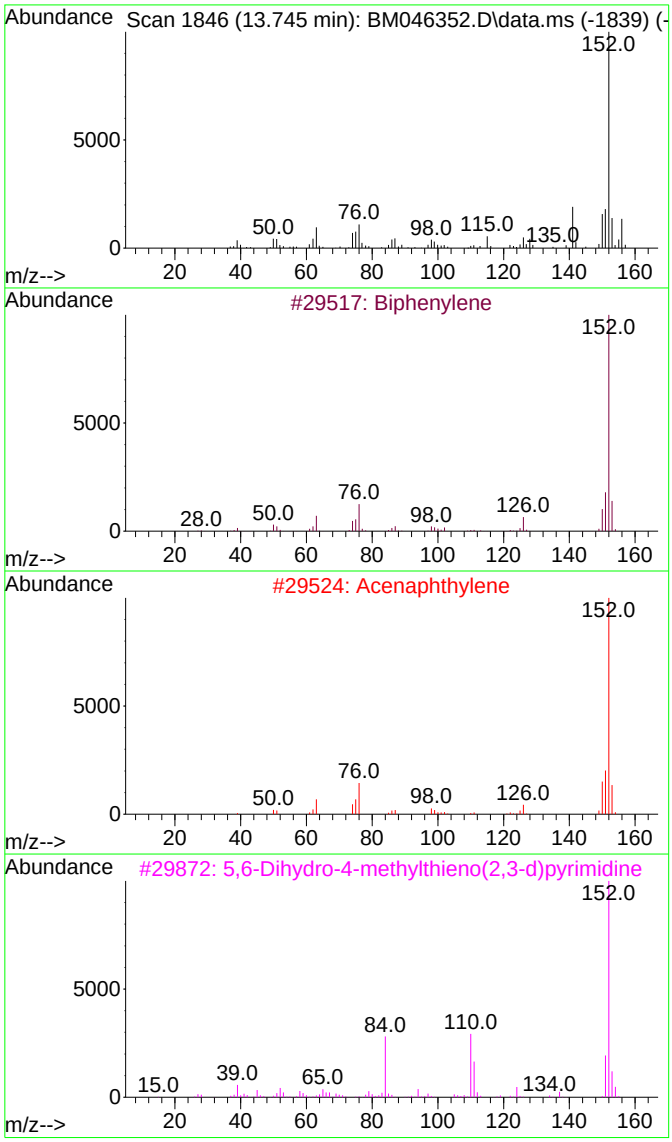
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

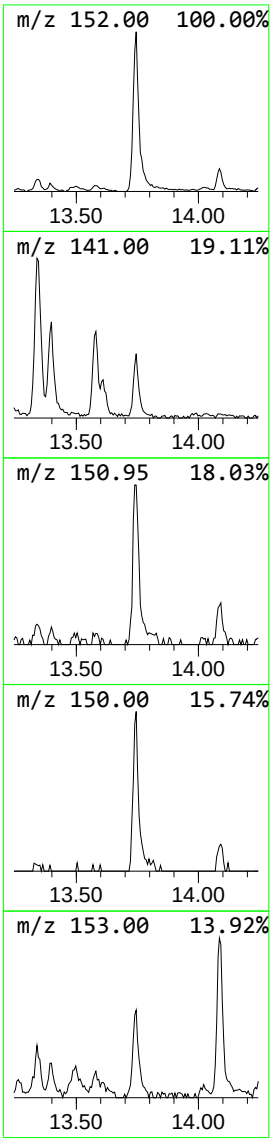
****

Peak Number 1 Biphenylene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.745	2.02 ng	163070	Acenaphthene-d10	14.021

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Biphenylene	152	C12H8	000259-79-0	76
2	Acenaphthylene	152	C12H8	000208-96-8	70
3	5,6-Dihydro-4-methylthieno(2,3-d...	152	C7H8N2S	092204-06-3	47
4	Benzaldehyde, 2-hydroxy-5-methoxy-	152	C8H8O3	000672-13-9	47
5	1-acenaphthenol, trifluoroacetat...	266	C14H9F3O2	1000376-45-2	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062824\
Data File : BM046352.D
Acq On : 28 Jun 2024 14:44
Operator : MA/JU
Sample : P3047-01 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EO-01-062424

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM062724.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

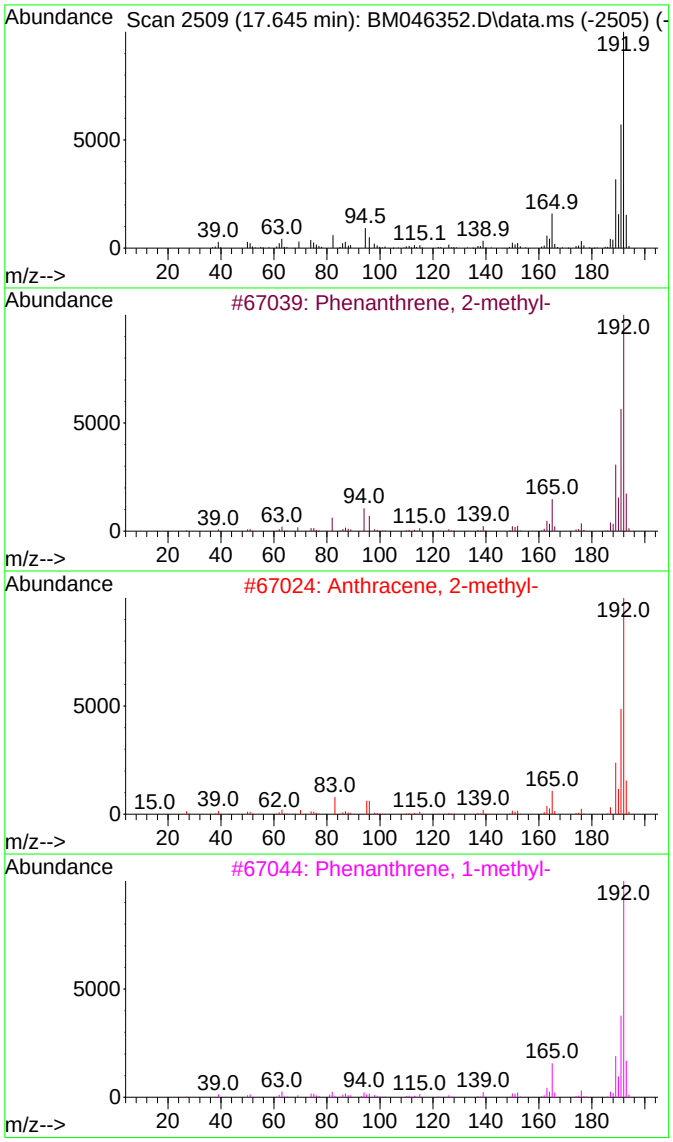
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

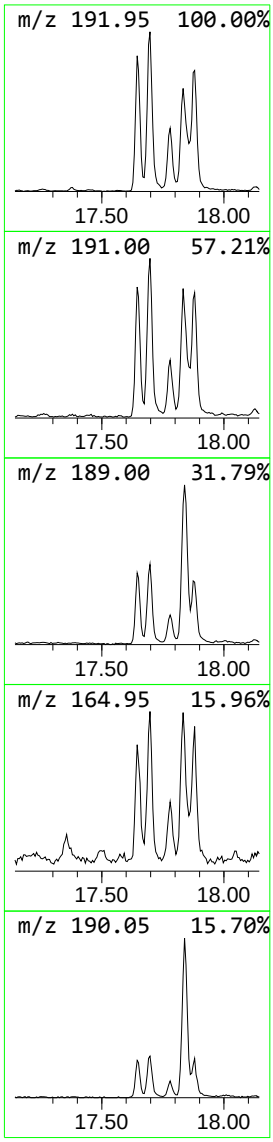
****

Peak Number 3 Phenanthrene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.645	3.30 ng	279817	Phenanthrene-d10	16.774

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
5	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	93





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Data File : BM046352.D
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Operator : MA/JU
Sample : P3047-01 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EO-01-062424

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM062724.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

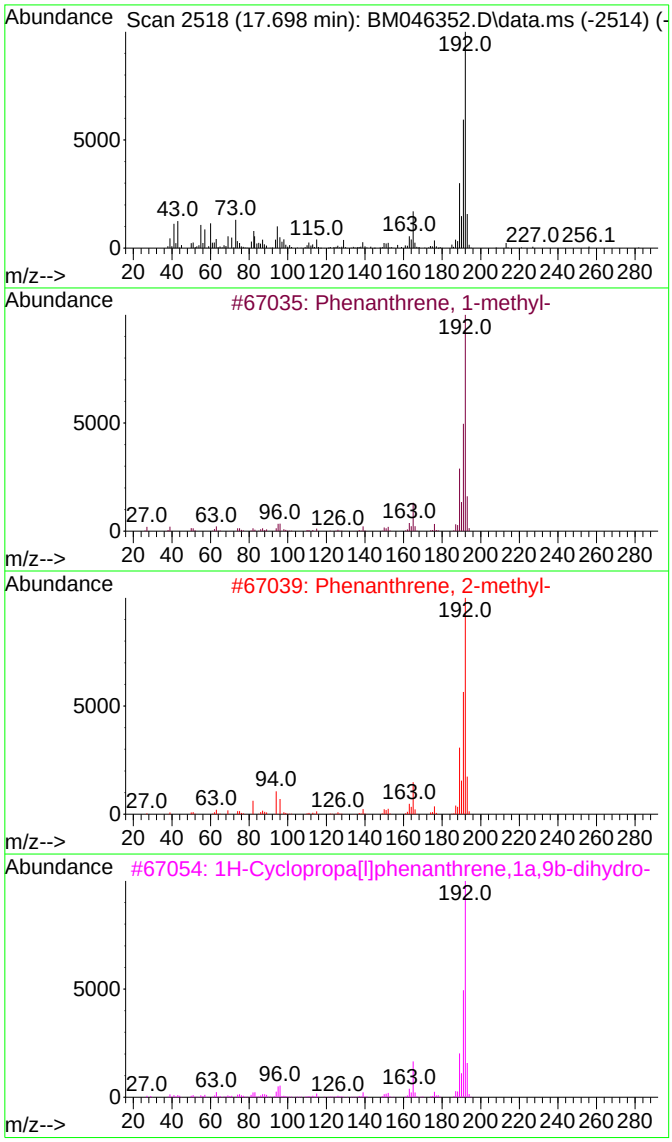
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

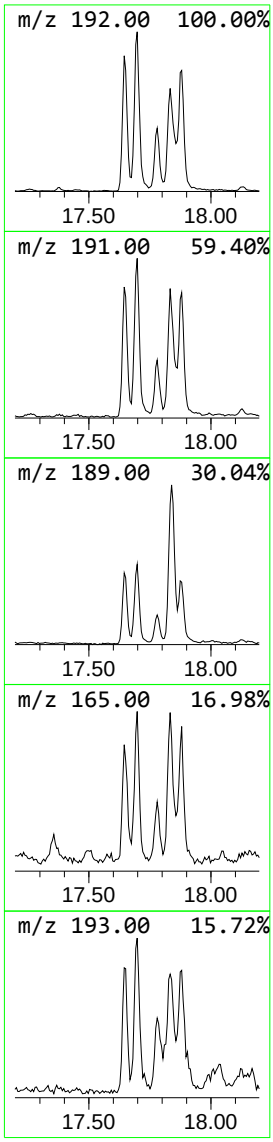
****

Peak Number 4 Phenanthrene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.698	6.01 ng	509897	Phenanthrene-d10	16.774

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
3	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	96
4	Anthracene, 2-methyl-	192	C15H12	000613-12-7	95
5	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	95





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Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
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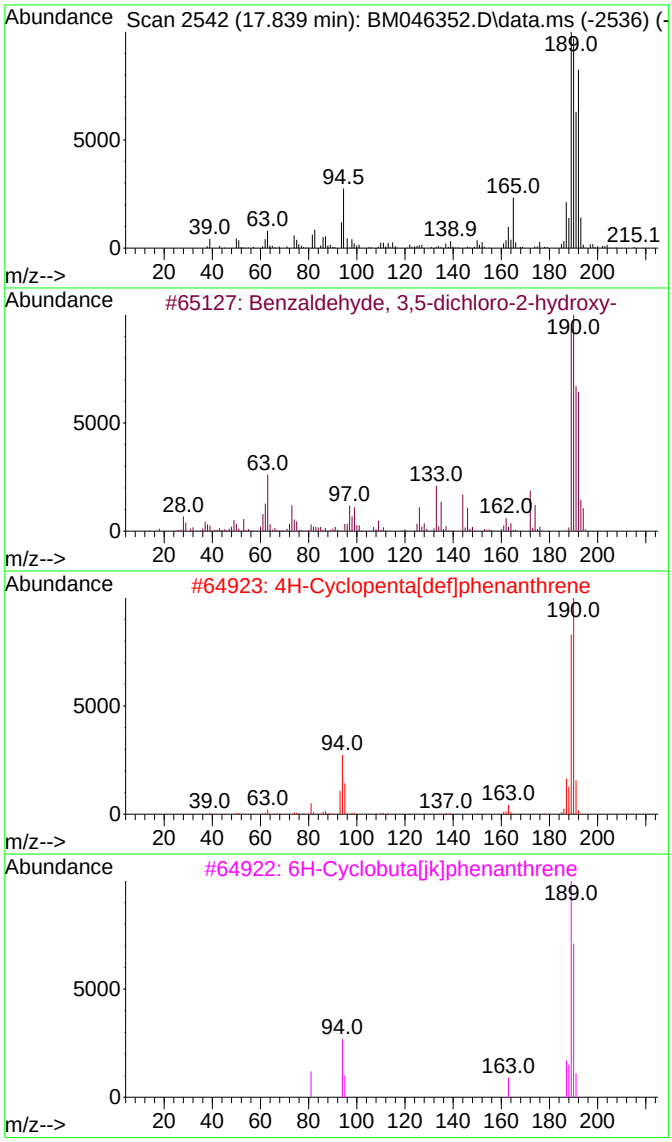
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM062724.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

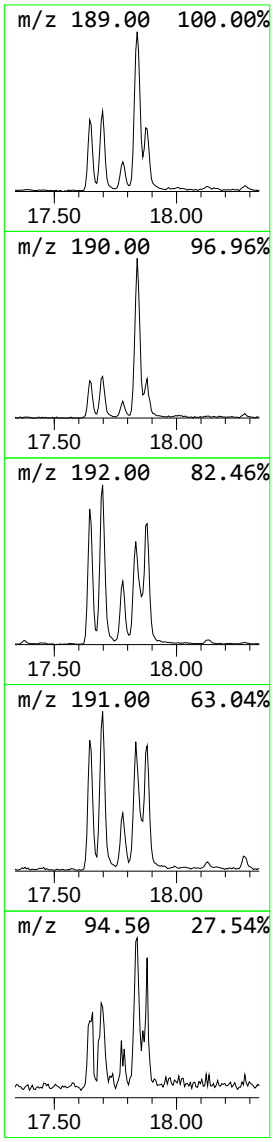
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzaldehyde, 3,5-dichloro-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.839	5.70 ng	483675	Phenanthrene-d10	16.774

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	64
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	60
3	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	49
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	46
5	Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	46





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Data File : BM046352.D
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Operator : MA/JU
Sample : P3047-01 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EO-01-062424

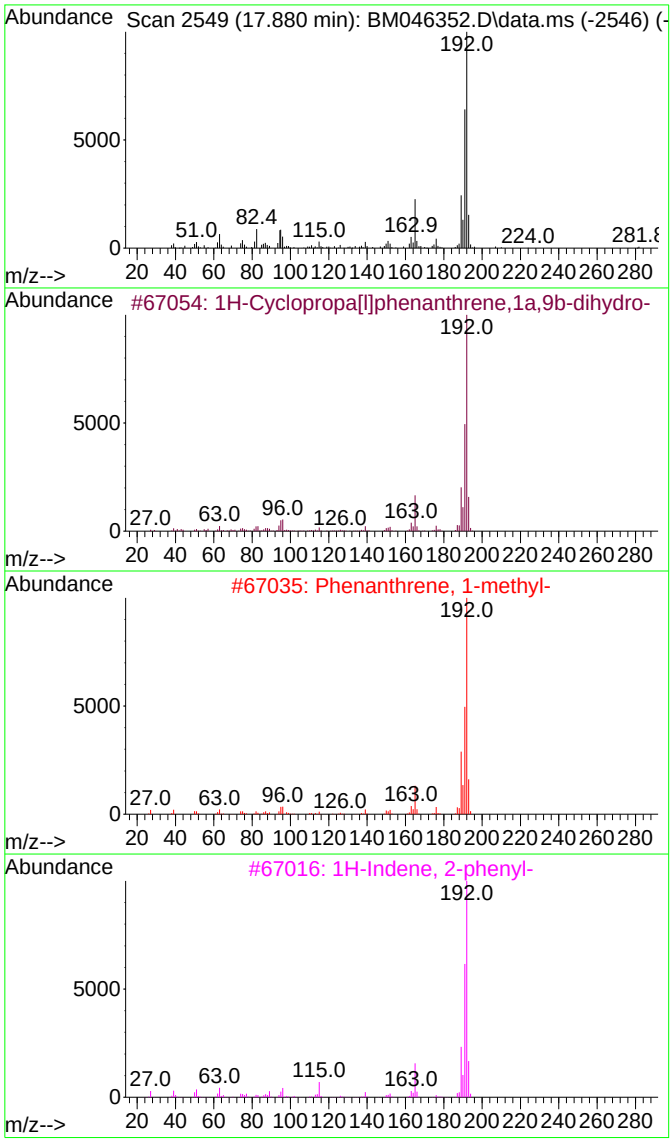
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM062724.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

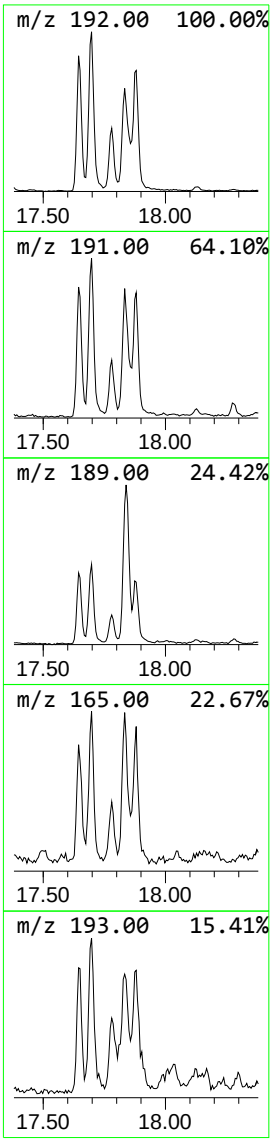
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 1H-Cyclopropa[l]phenanthren... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.880	3.39 ng	287364	Phenanthrene-d10	16.774

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	93
2	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
3	1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	93
4	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	91
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	91





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Sample : P3047-01 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM062724.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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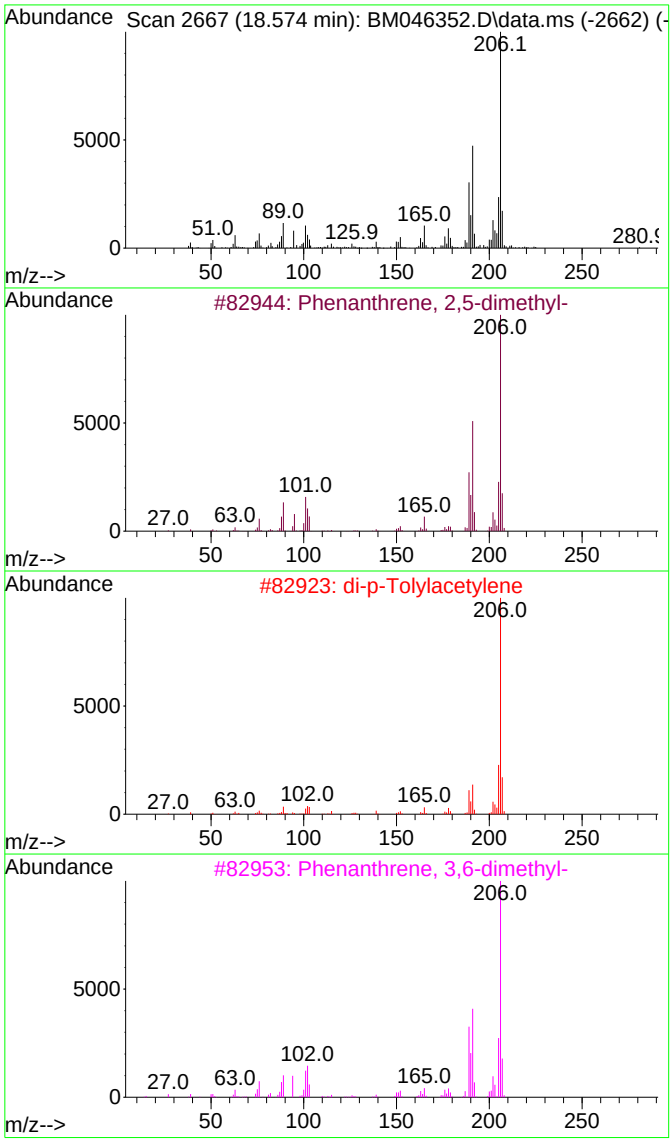
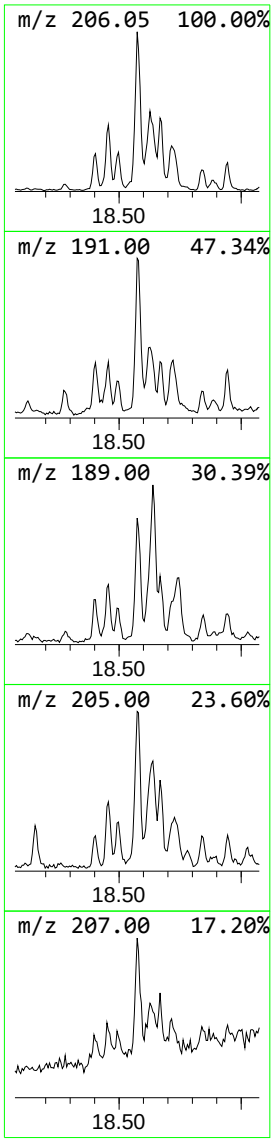
Peak Number 7 Phenanthrene, 2,5-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.574	3.60 ng	305684	Phenanthrene-d10	16.774

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	95
2	di-p-Tolylacetylene	206	C16H14	002789-88-0	93
3	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	93
4	Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	93
5	Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062824\
Data File : BM046352.D
Acq On : 28 Jun 2024 14:44
Operator : MA/JU
Sample : P3047-01 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EO-01-062424

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM062724.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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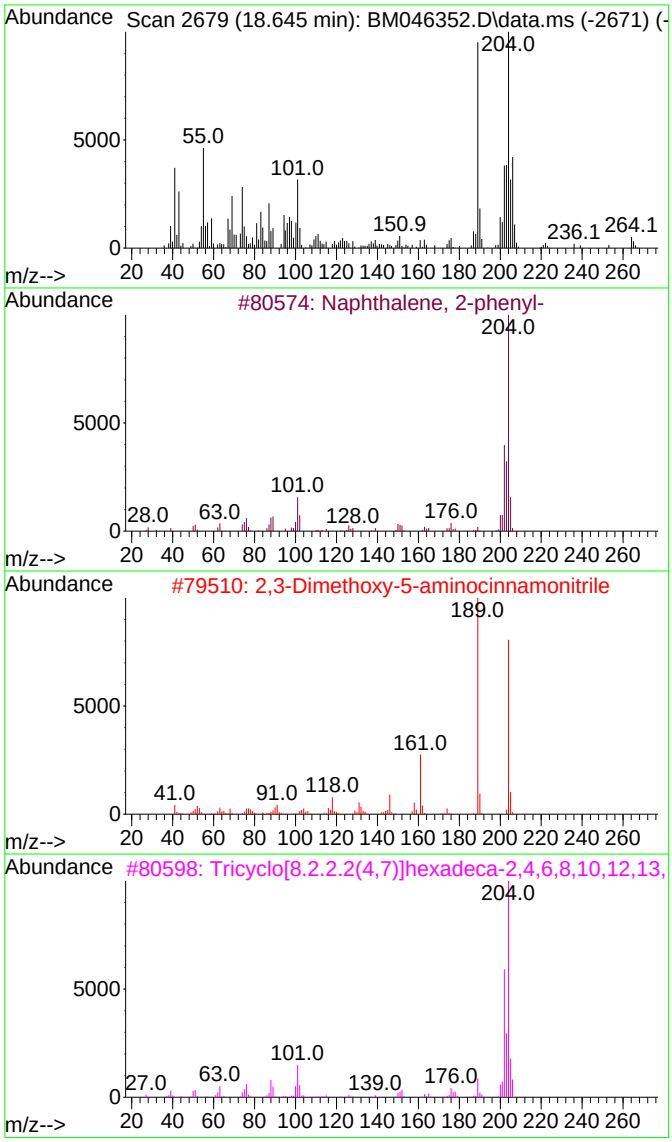
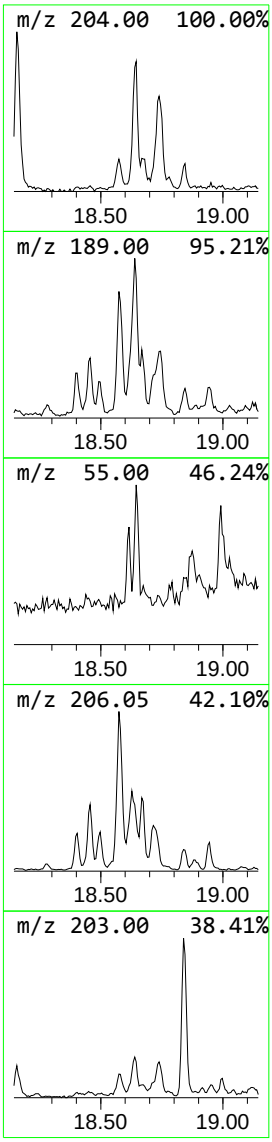
Peak Number 8 unknown18.645 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.645	4.24 ng	359527	Phenanthrene-d10	16.774

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	42
2	2,3-Dimethoxy-5-aminocinnamonitrile	204	C11H12N2O2	1000214-46-5	35
3	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	30
4	2-Chloro-4-phenylphenol	204	C12H9ClO	000092-04-6	27
5	Silane, dimethyl(2-methylpent-3-...	232	C12H28O2Si	1000347-64-8	22





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Data File : BM046352.D
Acq On : 28 Jun 2024 14:44
Operator : MA/JU
Sample : P3047-01 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EO-01-062424

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM062724.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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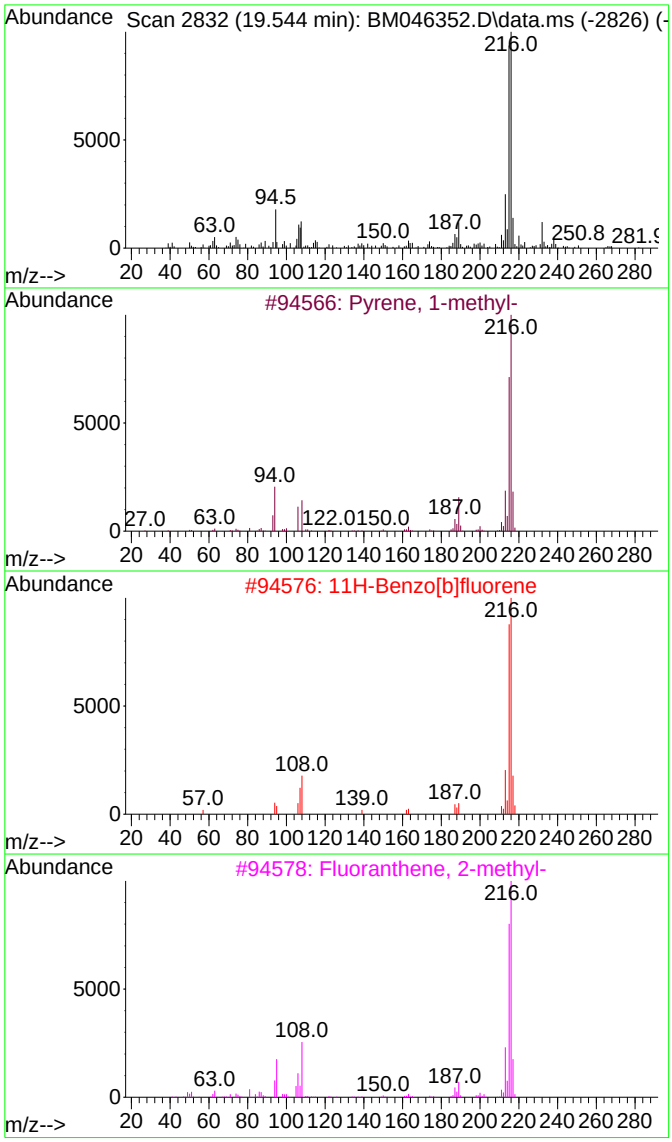
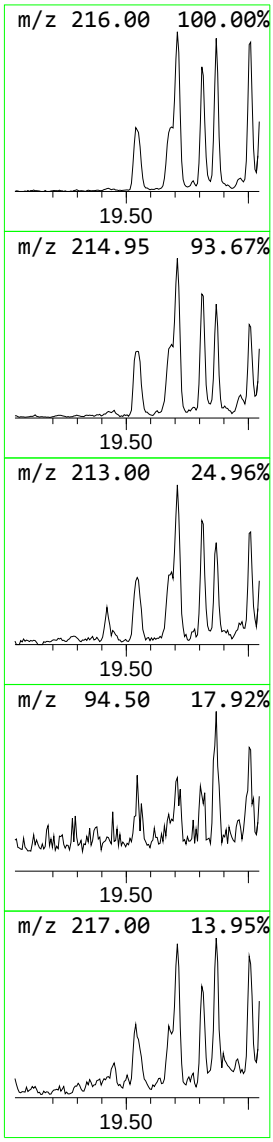
Peak Number 9 Pyrene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.544	2.92 ng	249340	Chrysene-d12	20.991

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	91
2	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
3	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	86
4	Pyrene, 2-methyl-	216	C17H12	003442-78-2	83
5	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	74





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Data File : BM046352.D
Acq On : 28 Jun 2024 14:44
Operator : MA/JU
Sample : P3047-01 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EO-01-062424

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM062724.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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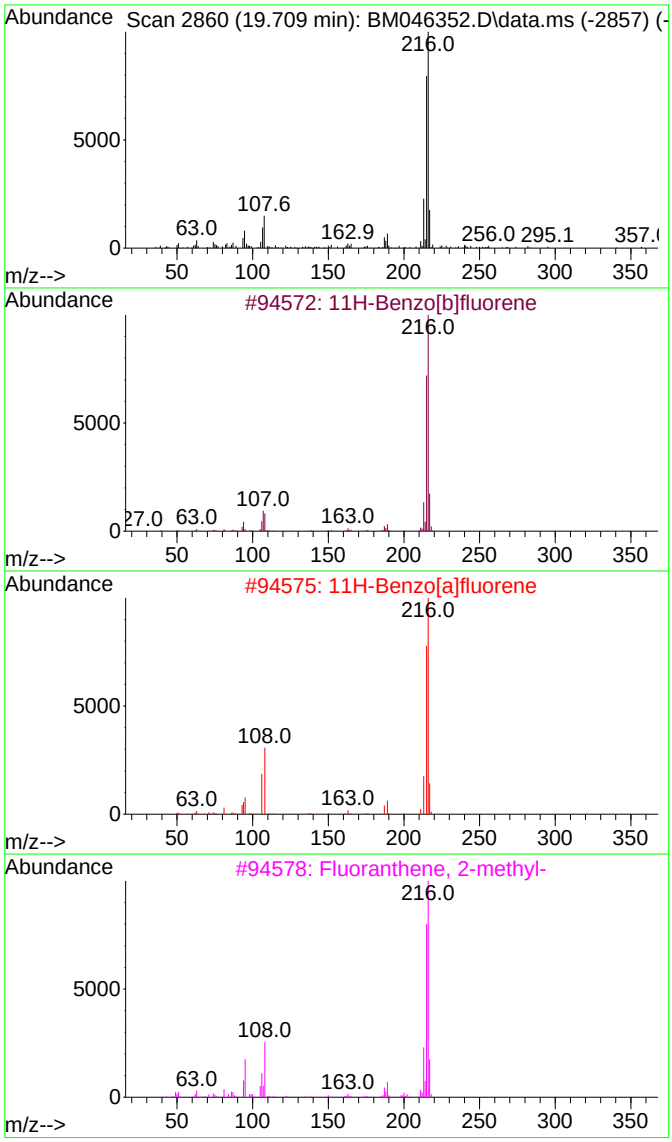
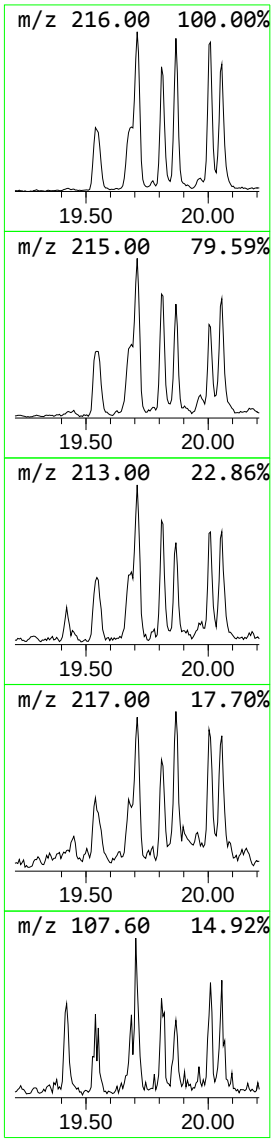
Peak Number 10 11H-Benzo[b]fluorene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.709	2.46 ng	210015	Chrysene-d12	20.991

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
2	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
3	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	81
4	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	81
5	Pyrene, 1-methyl-	216	C17H12	002381-21-7	59





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Data File : BM046352.D
Acq On : 28 Jun 2024 14:44
Operator : MA/JU
Sample : P3047-01 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EO-01-062424

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM062724.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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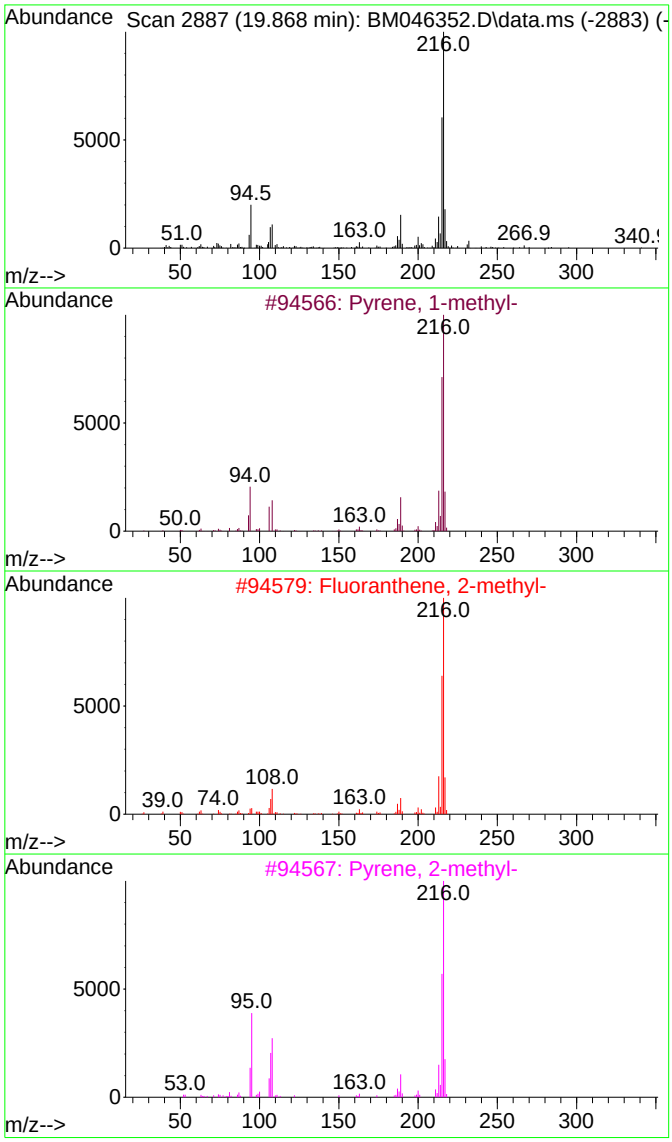
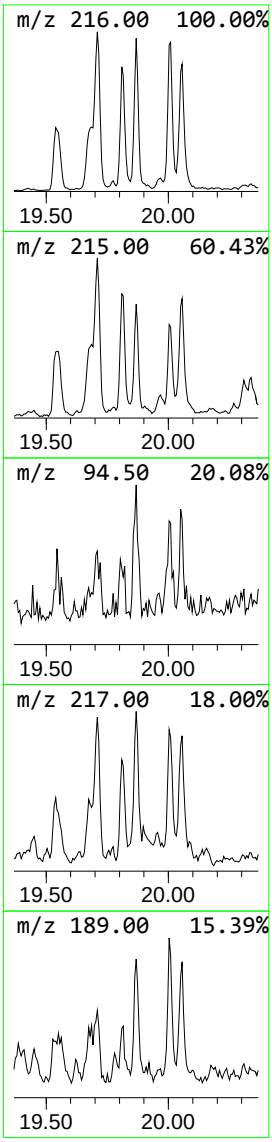
Peak Number 11 Fluoranthene, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.868	2.23 ng	190252	Chrysene-d12	20.991

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	98
2	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	96
3	Pyrene, 2-methyl-	216	C17H12	003442-78-2	94
4	Pyrene, 4-methyl-	216	C17H12	003353-12-6	93
5	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	83





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM062824\
Data File : BM046352.D
Acq On : 28 Jun 2024 14:44
Operator : MA/JU
Sample : P3047-01 10X
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EO-01-062424

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM062724.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
Biphenylene	13.745	2.0	ng	163070	3	14.021	1617560	20.0
Phenanthrene, 2...	17.645	3.3	ng	279817	4	16.774	1696760	20.0
Phenanthrene, 1...	17.698	6.0	ng	509897	4	16.774	1696760	20.0
Benzaldehyde, 3...	17.839	5.7	ng	483675	4	16.774	1696760	20.0
1H-Cyclopropa[l...	17.880	3.4	ng	287364	4	16.774	1696760	20.0
Phenanthrene, 2...	18.574	3.6	ng	305684	4	16.774	1696760	20.0
unknown18.645	18.645	4.2	ng	359527	4	16.774	1696760	20.0
Pyrene, 1-methyl-	19.544	2.9	ng	249340	5	20.991	1705850	20.0
11H-Benzo[b]flu...	19.709	2.5	ng	210015	5	20.991	1705850	20.0
Fluoranthene, 2...	19.868	2.2	ng	190252	5	20.991	1705850	20.0