

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110323\
 Data File : BM042579.D
 Acq On : 03 Nov 2023 20:23
 Operator : MA/JU
 Sample : 05220-02 5X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 Z-5(0-5)

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-BM103023.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM042579.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.452	378	385	391	rBV	137029	209965	11.91%	0.797%
2	7.081	656	662	673	rBV	123734	213485	12.11%	0.810%
3	7.916	795	804	814	rBV	219566	357276	20.27%	1.356%
4	8.445	886	894	900	rBV	24578	43213	2.45%	0.164%
5	9.116	1002	1008	1015	rBV	75176	121466	6.89%	0.461%
6	10.734	1277	1283	1288	rBV	274958	456470	25.90%	1.733%
7	10.786	1288	1292	1302	rVB	165747	271909	15.43%	1.032%
8	10.916	1307	1314	1319	rVB2	17359	29722	1.69%	0.113%
9	12.069	1504	1510	1516	rBV	22069	32438	1.84%	0.123%
10	12.286	1544	1547	1555	rVB2	17868	29982	1.70%	0.114%
11	12.392	1559	1565	1573	rVV	117147	179884	10.21%	0.683%
12	12.610	1593	1602	1609	rBV	241936	389341	22.09%	1.478%
13	13.186	1694	1700	1706	rBV	221735	305708	17.35%	1.160%
14	13.316	1718	1722	1731	rVB	28663	42023	2.38%	0.160%
15	13.404	1731	1737	1744	rVB	41230	66573	3.78%	0.253%
16	13.592	1765	1769	1772	rBV	29009	45984	2.61%	0.175%
17	13.722	1784	1791	1801	rBV2	80572	213588	12.12%	0.811%
18	13.804	1801	1805	1811	rBV	26915	40435	2.29%	0.153%
19	13.880	1811	1818	1823	rBV	153393	270618	15.36%	1.027%
20	13.933	1823	1827	1831	rBV	50157	76155	4.32%	0.289%
21	14.022	1837	1842	1850	rVB	49392	77372	4.39%	0.294%
22	14.116	1850	1858	1862	rBV	74153	117363	6.66%	0.445%
23	14.151	1862	1864	1868	rVB	28058	29953	1.70%	0.114%
24	14.298	1882	1889	1900	rBV	233572	409377	23.23%	1.554%
25	14.392	1900	1905	1910	rBV	31110	39677	2.25%	0.151%
26	14.569	1930	1935	1941	rVV	388725	552299	31.34%	2.096%
27	14.633	1941	1946	1952	rVB2	72491	111358	6.32%	0.423%
28	14.769	1962	1969	1975	rBV3	36094	83142	4.72%	0.316%
29	14.945	1992	1999	2007	rVB4	59393	140750	7.99%	0.534%
30	15.027	2007	2013	2018	rVB	45059	74748	4.24%	0.284%
31	15.163	2030	2036	2041	rBV2	66677	121909	6.92%	0.463%
32	15.221	2041	2046	2050	rVB2	46403	63286	3.59%	0.240%
33	15.339	2058	2066	2070	rBV3	58069	109033	6.19%	0.414%
34	15.374	2070	2072	2078	rVB2	23787	29018	1.65%	0.110%
35	15.439	2078	2083	2087	rBV	47339	72867	4.13%	0.277%
36	15.504	2087	2094	2099	rVB4	14894	35015	1.99%	0.133%

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Filtering: 5

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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-BM103023.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	15.557	2099	2103	2106	rBV	34698	47605	2.70%	0.181%
38	15.616	2107	2113	2124	rVB	188116	335465	19.04%	1.273%
39	15.733	2130	2133	2142	rVB5	33127	71538	4.06%	0.272%
40	15.898	2158	2161	2166	rVB2	33018	48596	2.76%	0.184%

41	15.957	2166	2171	2176	rVB3	22589	44369	2.52%	0.168%
42	16.063	2176	2189	2194	rBV3	172189	339875	19.29%	1.290%
43	16.210	2208	2214	2219	rBV2	63350	112031	6.36%	0.425%
44	16.263	2220	2223	2229	rVB4	25106	42581	2.42%	0.162%
45	16.610	2274	2282	2287	rBV2	70716	151691	8.61%	0.576%

46	16.663	2287	2291	2296	rBV2	50297	75500	4.28%	0.287%
47	16.763	2303	2308	2313	rBV	43317	76386	4.33%	0.290%
48	16.827	2313	2319	2322	rBV7	13956	29806	1.69%	0.113%
49	16.974	2339	2344	2345	rBV5	19433	32164	1.83%	0.122%
50	16.998	2345	2348	2351	rVV	58160	70680	4.01%	0.268%

51	17.039	2351	2355	2359	rVB3	38809	58329	3.31%	0.221%
52	17.139	2367	2372	2380	rVB	150771	216839	12.30%	0.823%
53	17.321	2397	2403	2406	rBV	466678	659558	37.43%	2.504%
54	17.363	2406	2410	2419	rVB	937693	1249321	70.89%	4.742%
55	17.457	2420	2426	2431	rBV	275585	425018	24.12%	1.613%

56	17.739	2470	2474	2479	rVB	90724	110092	6.25%	0.418%
57	17.833	2486	2490	2496	rBV	44820	64882	3.68%	0.246%
58	17.898	2496	2501	2510	rVB	118391	189318	10.74%	0.719%
59	18.045	2520	2526	2531	rVB	76083	156968	8.91%	0.596%
60	18.186	2542	2550	2555	rVV2	249762	536851	30.46%	2.038%

61	18.239	2555	2559	2565	rVV	257046	377284	21.41%	1.432%
62	18.321	2569	2573	2577	rVV	114408	160907	9.13%	0.611%
63	18.392	2578	2585	2588	rVV2	329727	636265	36.11%	2.415%
64	18.427	2588	2591	2597	rVB2	185759	273672	15.53%	1.039%
65	18.580	2615	2617	2622	rVB	40044	46921	2.66%	0.178%

66	18.692	2630	2636	2641	rBV2	112966	219631	12.46%	0.834%
67	18.845	2659	2662	2668	rBV3	31877	63861	3.62%	0.242%
68	18.904	2668	2672	2674	rBV2	39316	50753	2.88%	0.193%
69	18.980	2682	2685	2688	rVB	60079	69986	3.97%	0.266%
70	19.021	2688	2692	2696	rVB4	65971	83011	4.71%	0.315%

71	19.068	2697	2700	2702	rBV	68595	80918	4.59%	0.307%
72	19.098	2702	2705	2710	rVB	152081	205686	11.67%	0.781%
73	19.239	2726	2729	2730	rBV2	35589	39301	2.23%	0.149%
74	19.380	2747	2753	2759	rBV	845468	1244680	70.63%	4.724%
75	19.527	2774	2778	2783	rVB	159514	201488	11.43%	0.765%

76	19.627	2792	2795	2797	rBV	67904	91768	5.21%	0.348%
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77	19.704	2804	2808	2810	rBV	79285	126516	7.18%	0.480%
78	19.745	2810	2815	2821	rVV	1189371	1550629	87.99%	5.886%
79	19.804	2821	2825	2830	rVB3	57185	93389	5.30%	0.354%
80	19.933	2842	2847	2851	rBV	361881	460604	26.14%	1.748%
81	20.068	2864	2870	2876	rVB3	100617	217921	12.37%	0.827%
82	20.192	2887	2891	2894	rBV5	107996	201597	11.44%	0.765%
83	20.233	2895	2898	2904	rVB	227460	310627	17.63%	1.179%
84	20.333	2911	2915	2919	rBV	205240	234275	13.29%	0.889%
85	20.392	2921	2925	2928	rVV	134548	157969	8.96%	0.600%
86	20.527	2945	2948	2952	rBV	140795	183592	10.42%	0.697%
87	20.574	2953	2956	2963	rVB	159347	226059	12.83%	0.858%
88	21.145	3050	3053	3057	rBV	211883	275068	15.61%	1.044%
89	21.227	3064	3067	3070	rBV2	84214	118375	6.72%	0.449%
90	21.298	3076	3079	3085	rVB2	149083	196081	11.13%	0.744%
91	21.503	3107	3114	3118	rBV3	807419	1762242	100.00%	6.689%
92	21.539	3118	3120	3125	rVB	528732	628406	35.66%	2.385%
93	21.833	3167	3170	3174	rBV	244712	302169	17.15%	1.147%
94	22.403	3263	3267	3274	rVB2	443312	666707	37.83%	2.531%
95	23.033	3370	3374	3380	rVB	414091	619020	35.13%	2.350%
96	23.180	3395	3399	3413	rVB	282429	885739	50.26%	3.362%
97	23.750	3493	3496	3502	rVB	386834	642419	36.45%	2.438%
98	23.815	3503	3507	3514	rVB	354117	650451	36.91%	2.469%
99	23.927	3521	3526	3531	rBV	322308	571321	32.42%	2.169%
100	24.603	3637	3641	3659	rVB3	318457	809194	45.92%	3.071%

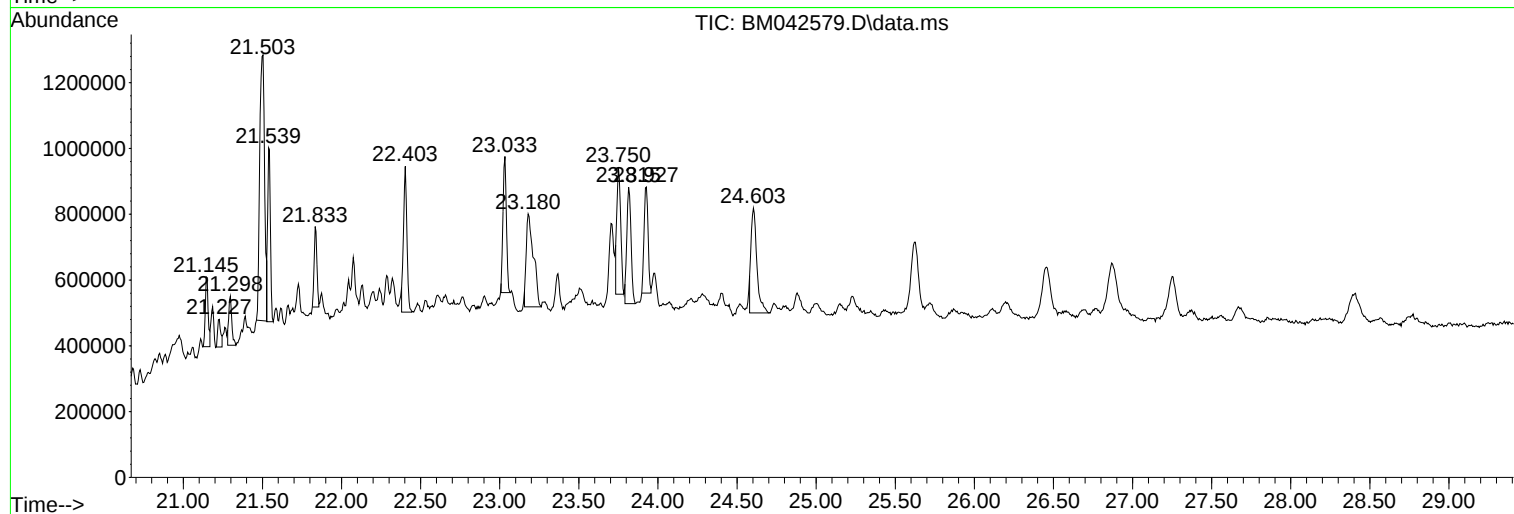
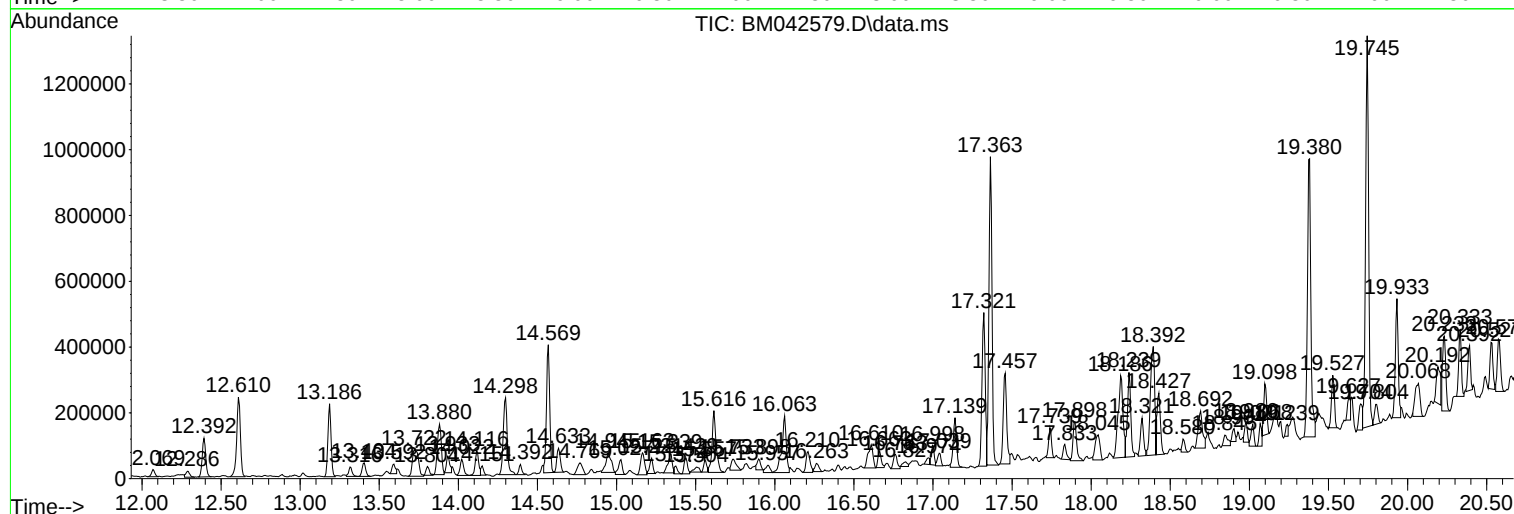
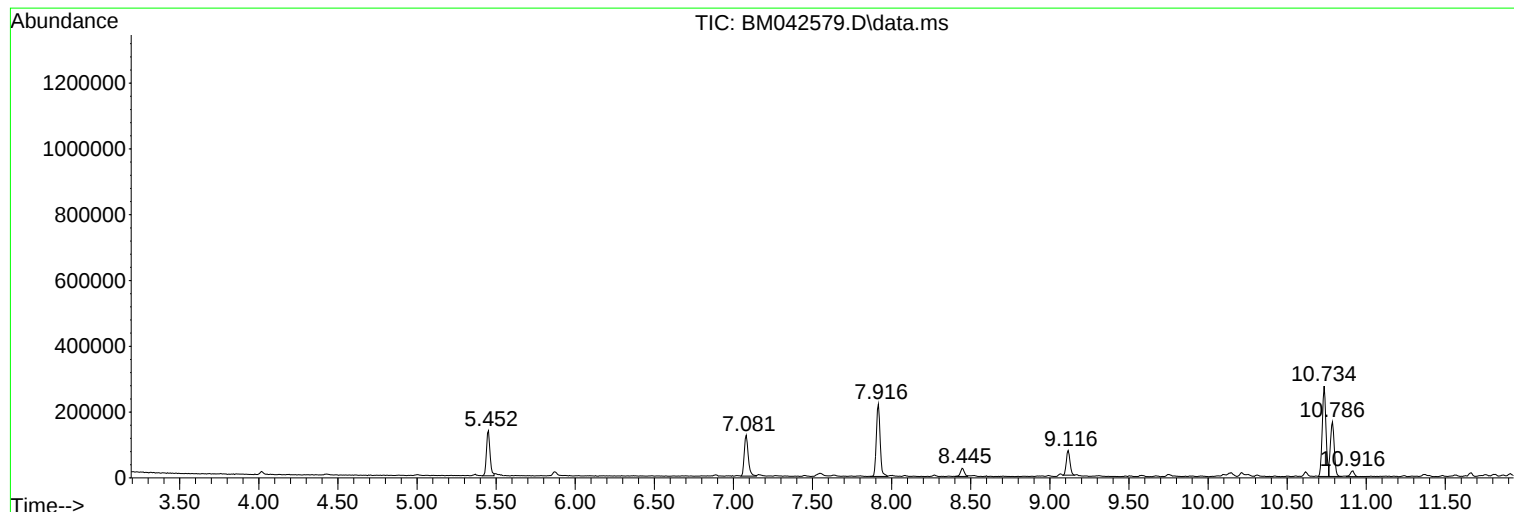
Sum of corrected areas: 26345367

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Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
Z-5(0-5)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM103023.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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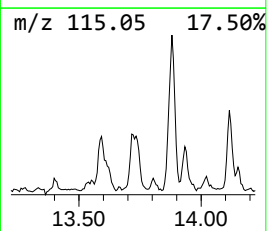
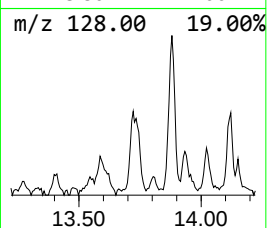
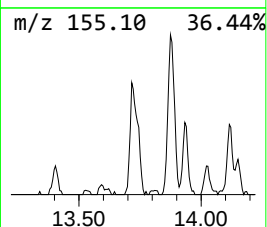
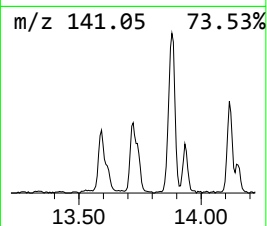
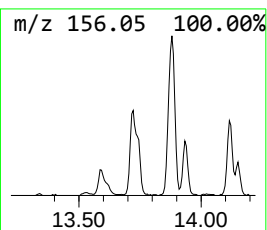
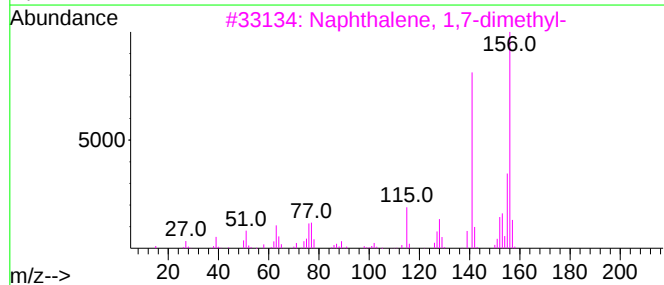
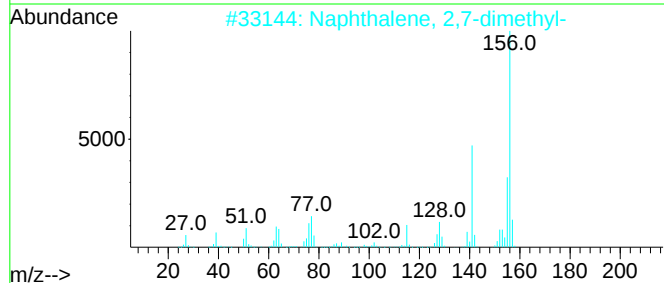
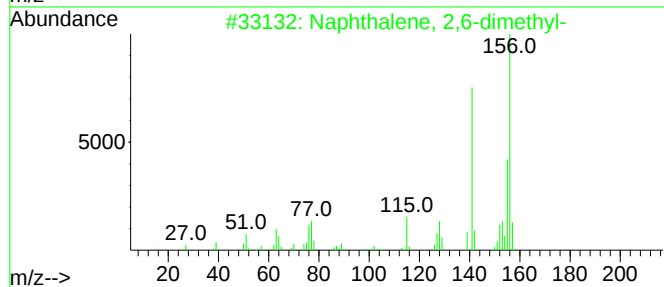
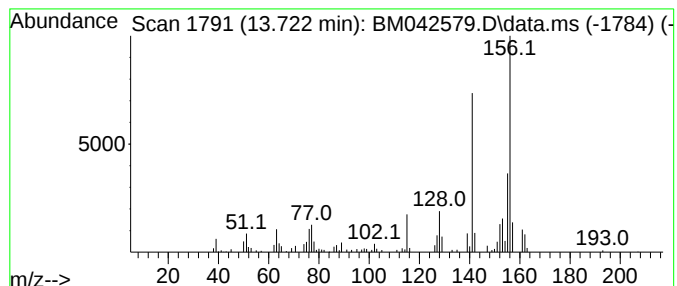
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM103023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Naphthalene, 2,6-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.722	7.73 ng	213588	Acenaphthene-d10	14.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
2			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
4			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
5			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95



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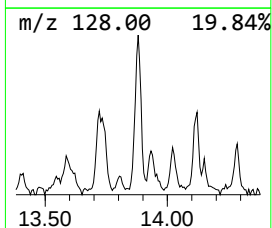
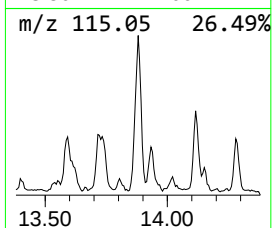
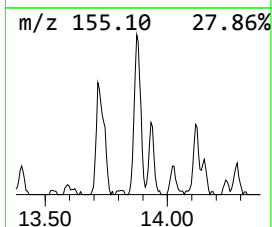
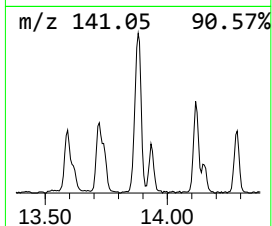
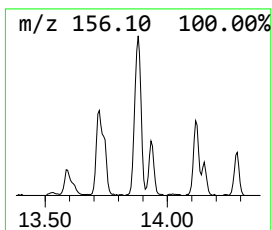
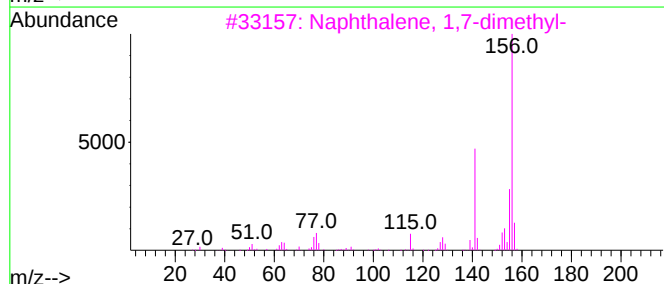
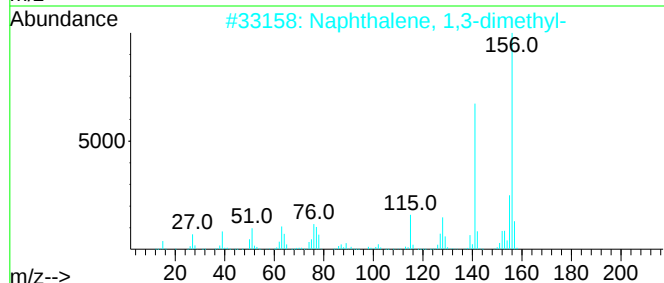
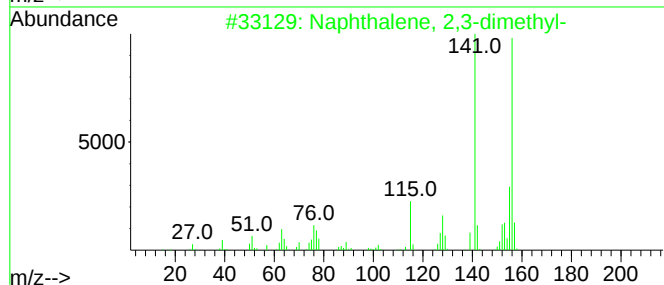
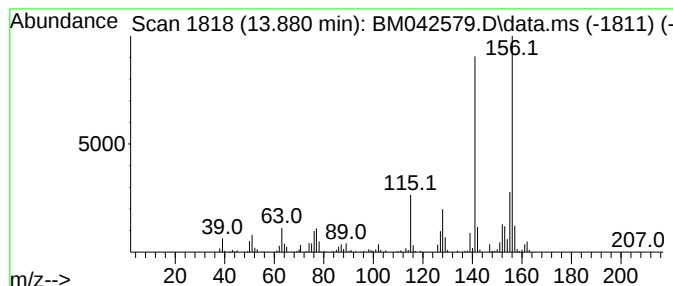
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM103023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Naphthalene, 2,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.880	9.80 ng	270618	Acenaphthene-d10	14.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
3			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
4			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
5			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95



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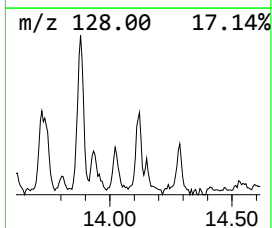
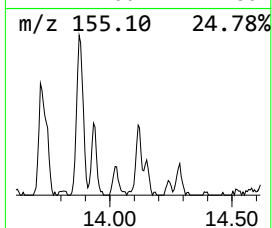
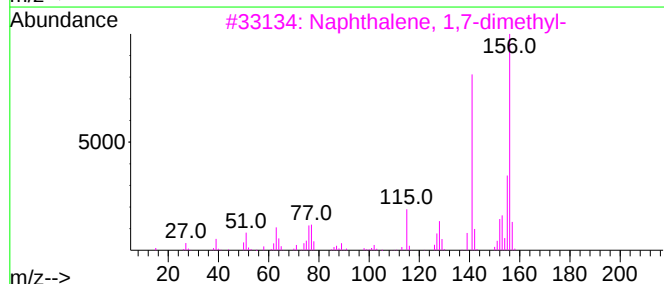
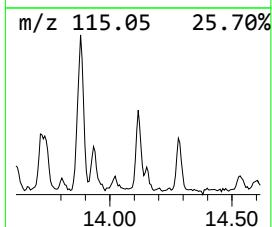
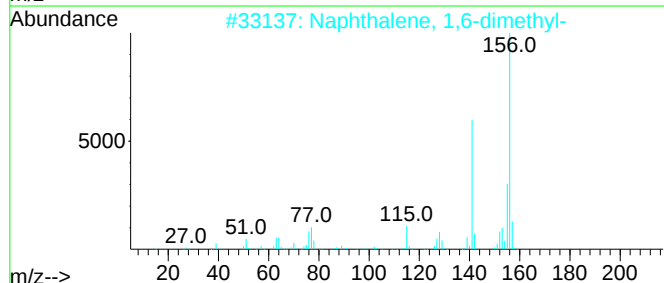
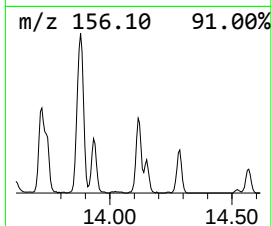
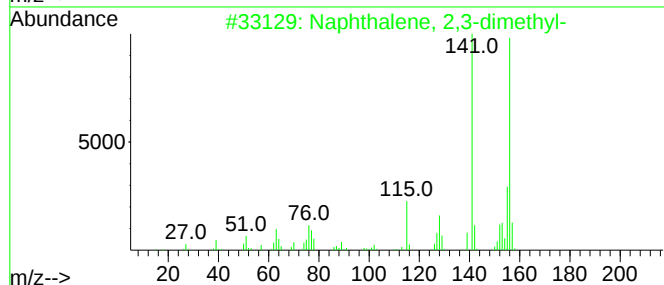
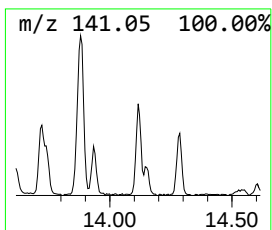
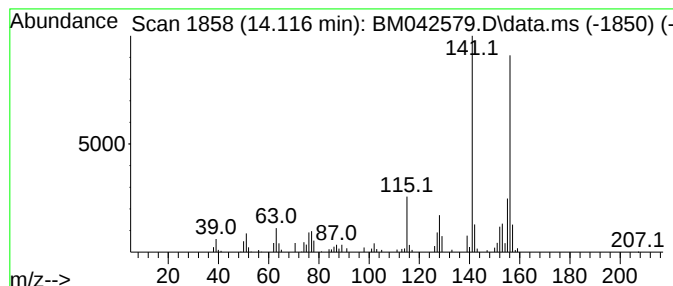
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Naphthalene, 1,6-dimethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.116	4.25 ng	117363	Acenaphthene-d10	14.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
3			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
4			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
5			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110323\
Data File : BM042579.D
Acq On : 03 Nov 2023 20:23
Operator : MA/JU
Sample : 05220-02 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

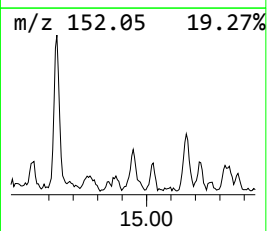
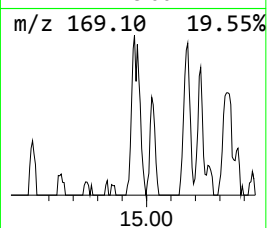
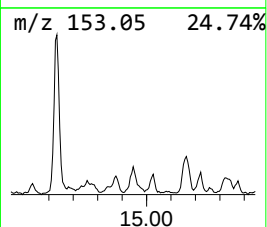
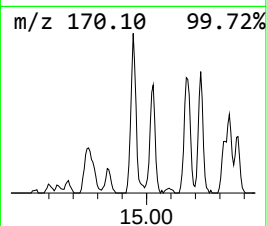
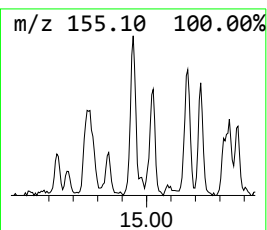
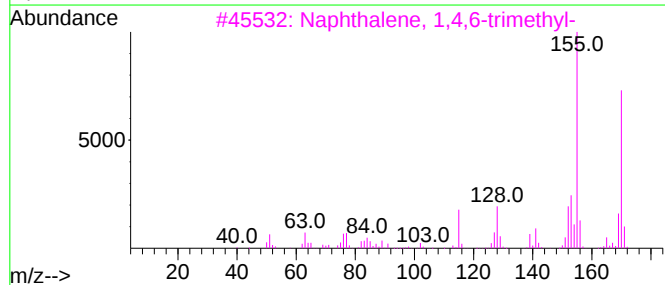
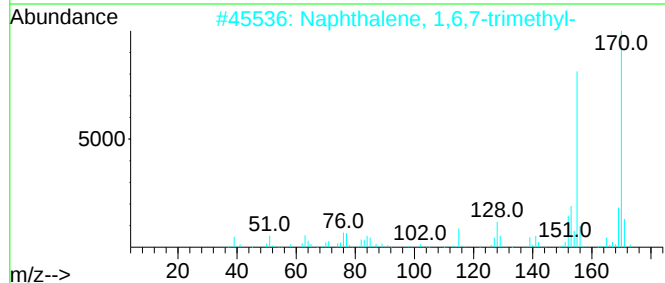
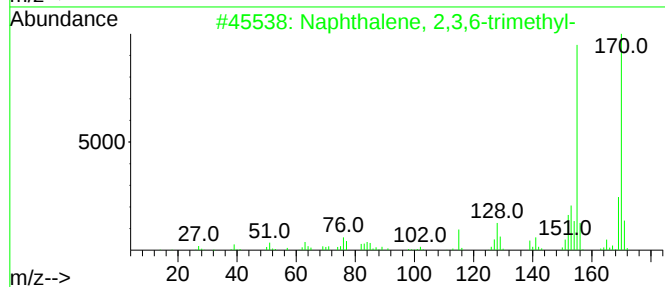
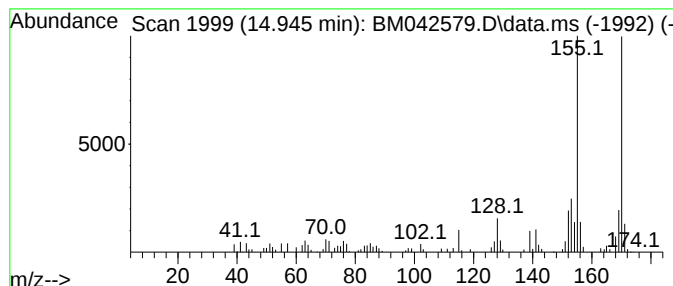
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Naphthalene, 2,3,6-trimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.945	5.10 ng	140750	Acenaphthene-d10	14.569	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	97
2	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	96
3	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	95
4	Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	93
5	3-(2-Methyl-propenyl)-1H-indene	170	C13H14	1000187-78-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110323\
Data File : BM042579.D
Acq On : 03 Nov 2023 20:23
Operator : MA/JU
Sample : 05220-02 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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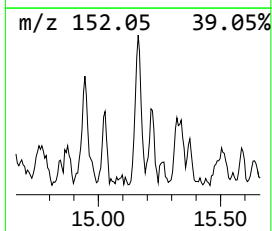
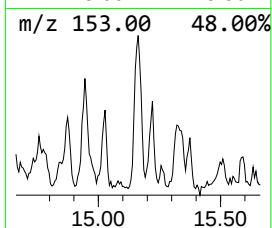
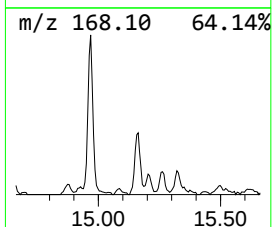
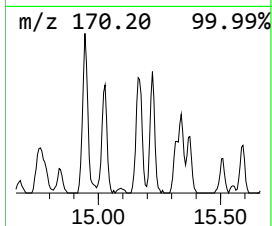
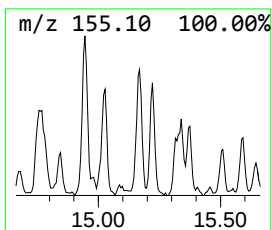
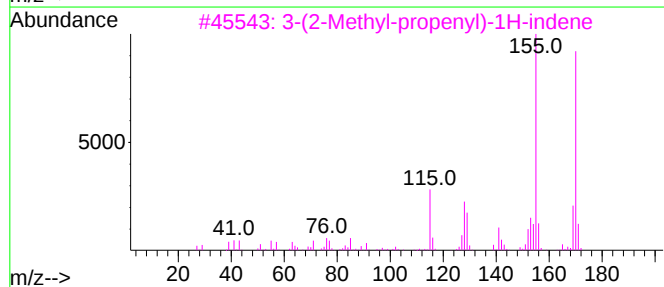
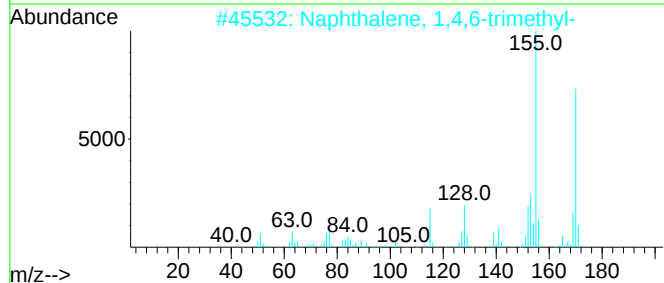
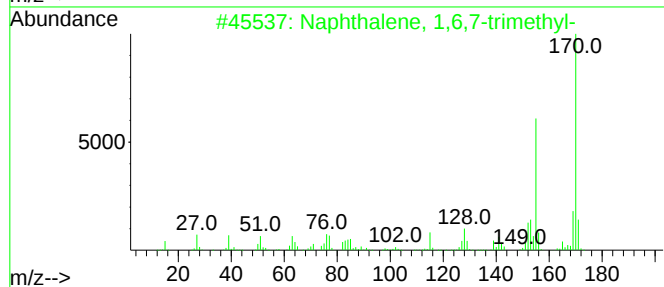
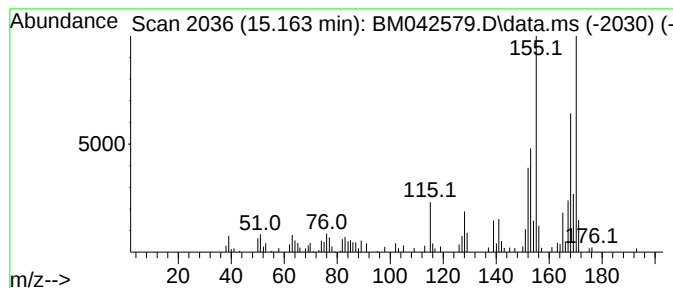
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Naphthalene, 1,6,7-trimethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.163	4.41 ng	121909	Acenaphthene-d10	14.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	87
2			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	83
3			3-(2-Methyl-propenyl)-1H-indene	170	C13H14	1000187-78-5	70
4			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	64
5			Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110323\
Data File : BM042579.D
Acq On : 03 Nov 2023 20:23
Operator : MA/JU
Sample : 05220-02 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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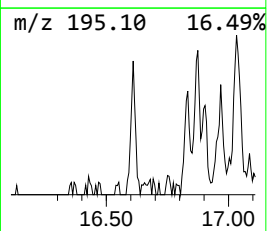
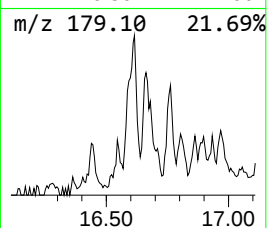
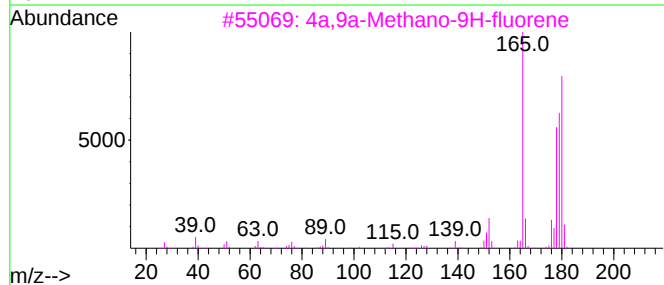
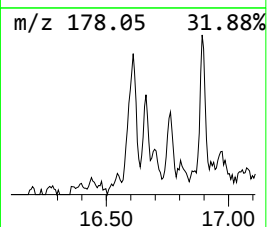
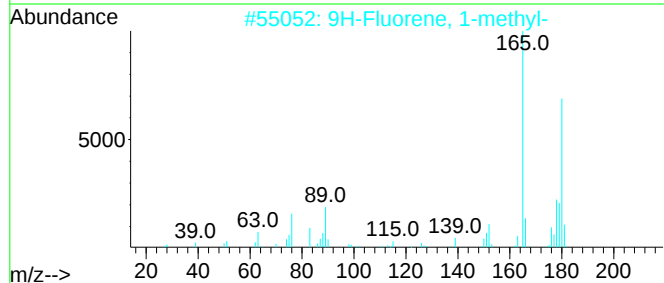
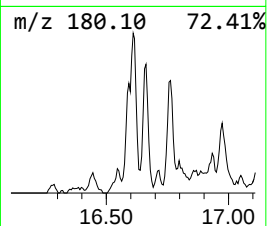
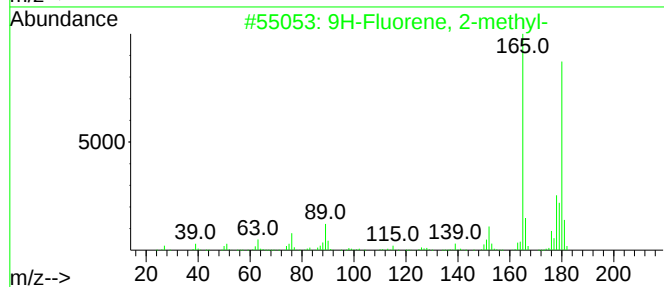
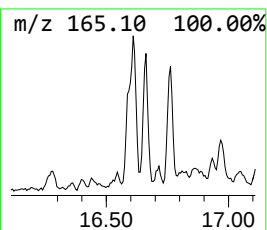
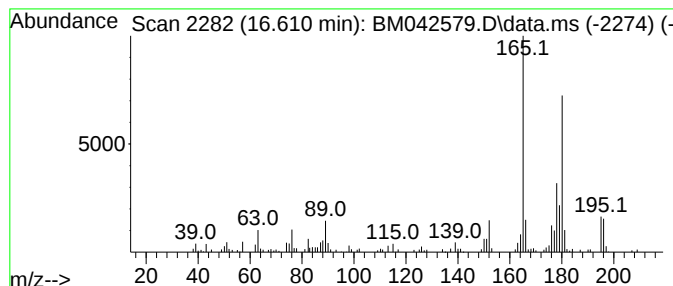
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 9H-Fluorene, 2-methyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.610	4.60 ng	151691	Phenanthrene-d10	17.321

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		9H-Fluorene, 2-methyl-	180	C14H12	001430-97-3	96
2		9H-Fluorene, 1-methyl-	180	C14H12	001730-37-6	95
3		4a,9a-Methano-9H-fluorene	180	C14H12	019540-84-2	91
4		9H-Fluorene, 9-methyl-	180	C14H12	002523-37-7	89
5		9H-Fluorene, 3-methyl-	180	C14H12	002523-39-9	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110323\
Data File : BM042579.D
Acq On : 03 Nov 2023 20:23
Operator : MA/JU
Sample : 05220-02 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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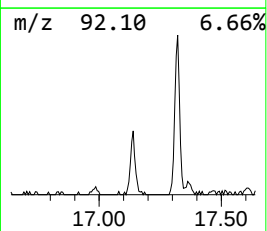
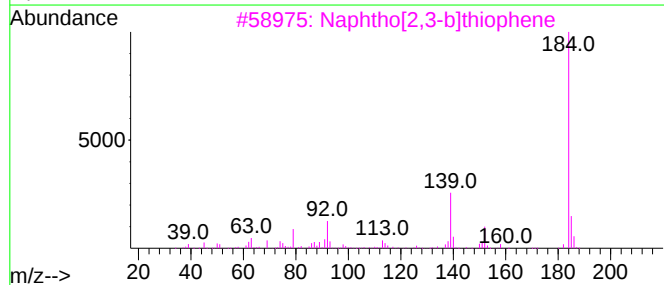
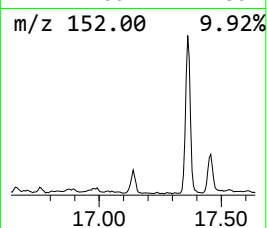
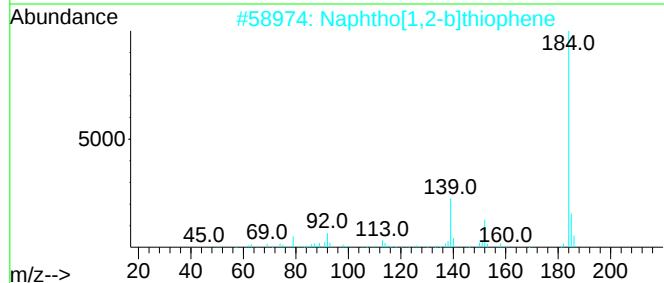
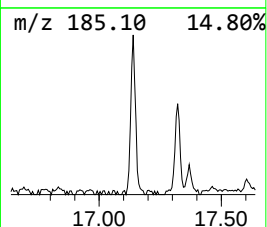
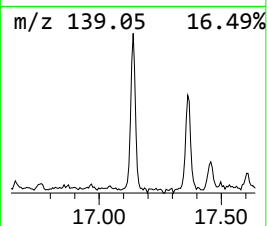
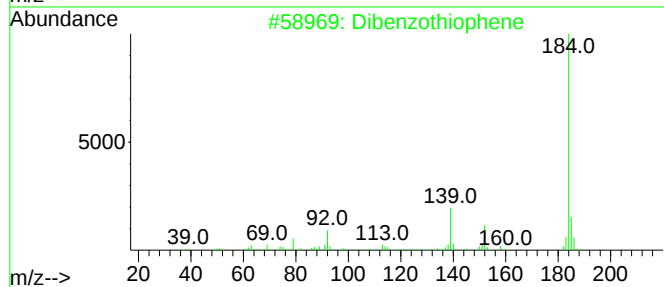
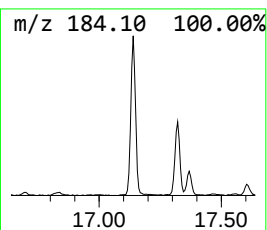
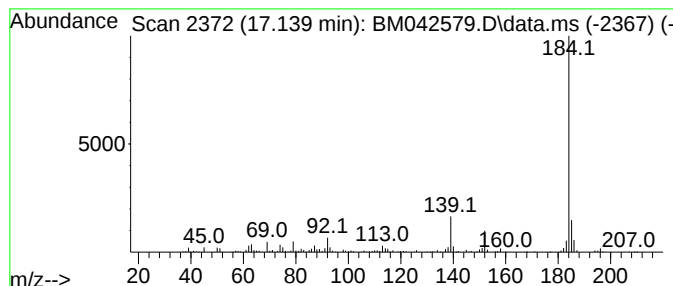
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Dibenzothiophene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.139	6.58 ng	216839	Phenanthrene-d10	17.321

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzothiophene	184	C12H8S	000132-65-0	97
2			Naphtho[1,2-b]thiophene	184	C12H8S	000234-41-3	94
3			Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	94
4			Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	94
5			Azuleno(2,1-b)thiophene	184	C12H8S	000248-13-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110323\
Data File : BM042579.D
Acq On : 03 Nov 2023 20:23
Operator : MA/JU
Sample : 05220-02 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

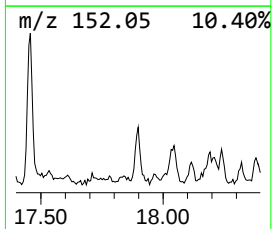
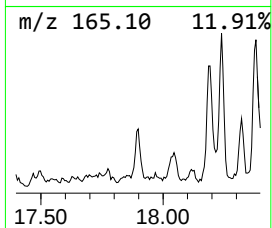
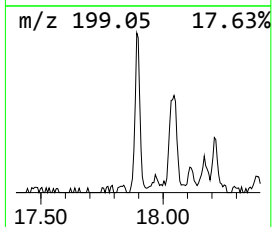
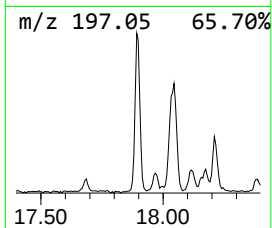
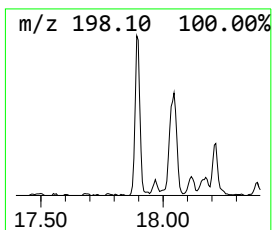
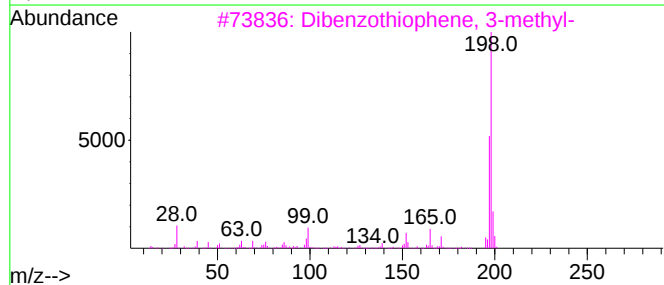
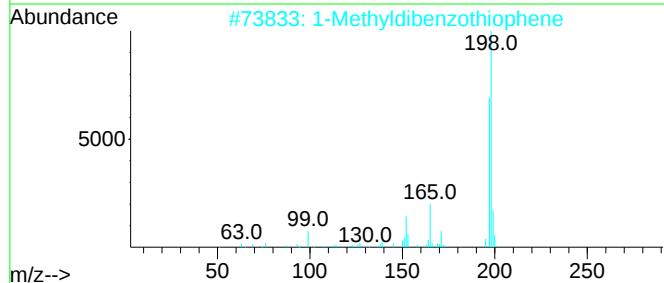
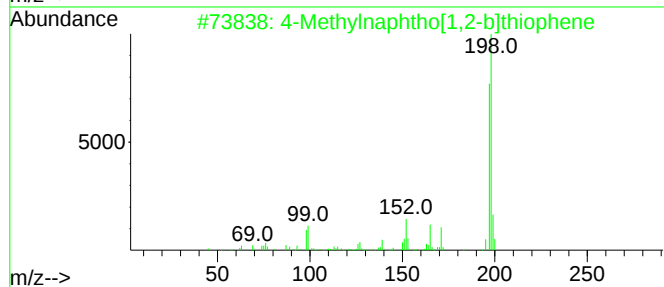
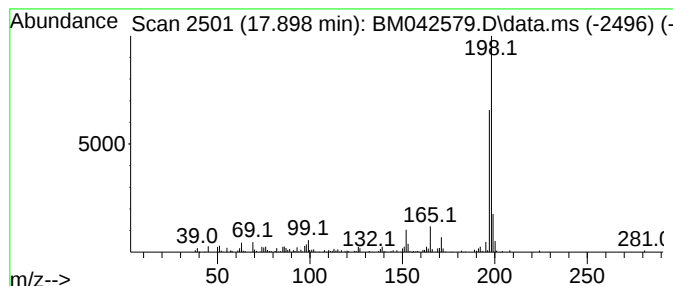
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 4-Methylnaphtho[1,2-b]thiop... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.898	5.74 ng	189318	Phenanthrene-d10	17.321	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Methylnaphtho[1,2-b]thiophene	198	C13H10S	067388-11-8	97
2	1-Methyldibenzothiophene	198	C13H10S	031317-07-4	95
3	Dibenzothiophene, 3-methyl-	198	C13H10S	016587-52-3	94
4	Dibenzothiophene, 4-methyl-	198	C13H10S	007372-88-5	94
5	2-Methyldibenzothiophene	198	C13H10S	1000383-55-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110323\
Data File : BM042579.D
Acq On : 03 Nov 2023 20:23
Operator : MA/JU
Sample : 05220-02 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

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ClientSampleId :
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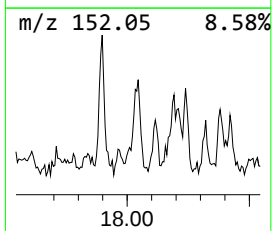
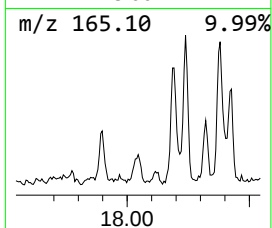
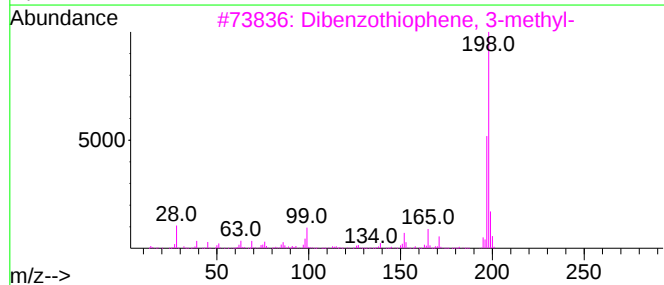
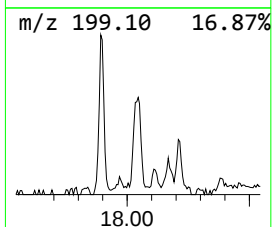
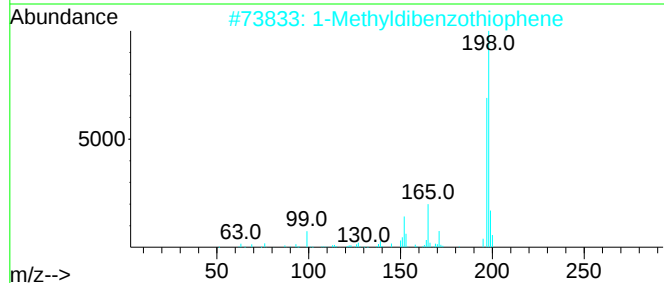
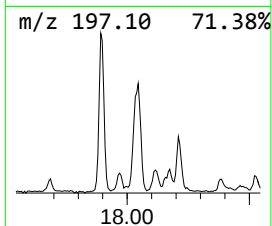
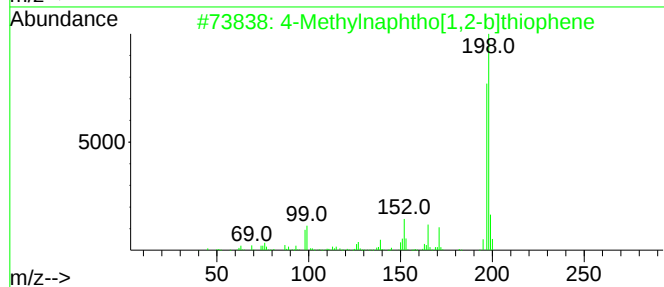
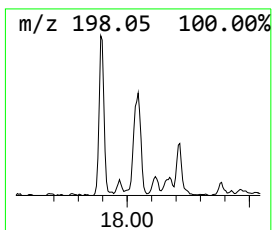
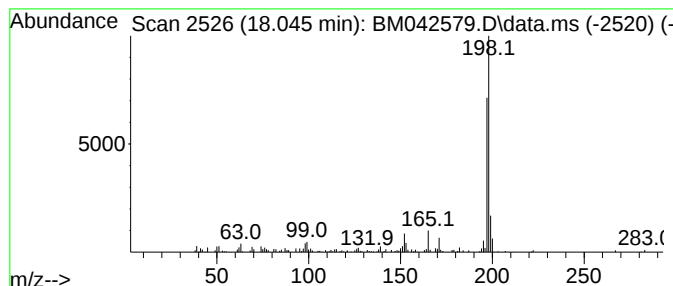
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 1-Methyldibenzothiophene Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.045	4.76 ng	156968	Phenanthrene-d10	17.321

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Methylnaphtho[1,2-b]thiophene	198	C13H10S	067388-11-8	96
2			1-Methyldibenzothiophene	198	C13H10S	031317-07-4	95
3			Dibenzothiophene, 3-methyl-	198	C13H10S	016587-52-3	94
4			Dibenzothiophene, 4-methyl-	198	C13H10S	007372-88-5	94
5			2-Methyldibenzothiophene	198	C13H10S	1000383-55-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110323\
Data File : BM042579.D
Acq On : 03 Nov 2023 20:23
Operator : MA/JU
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Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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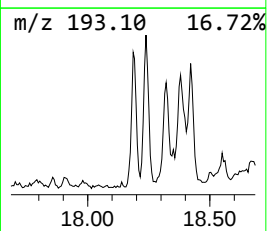
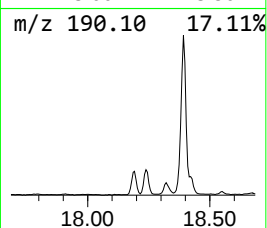
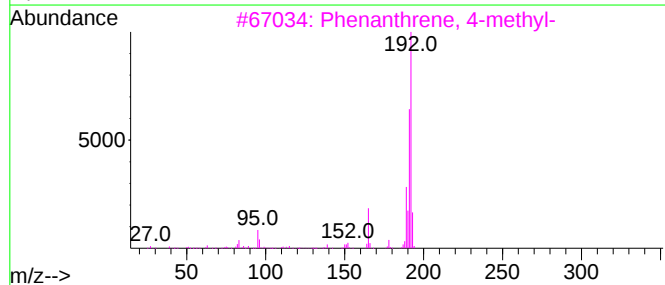
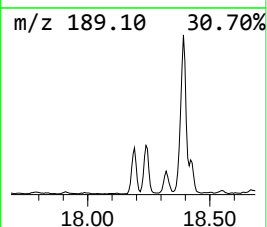
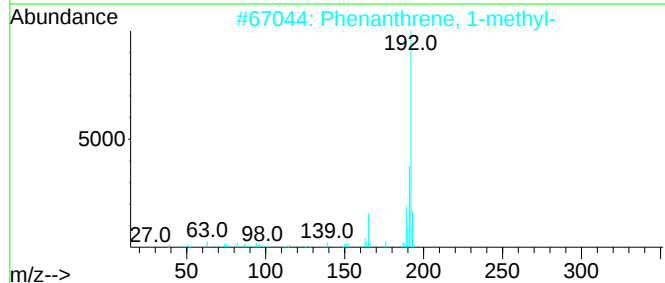
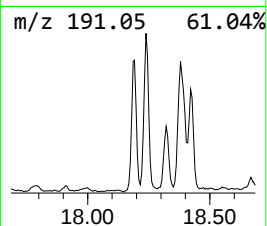
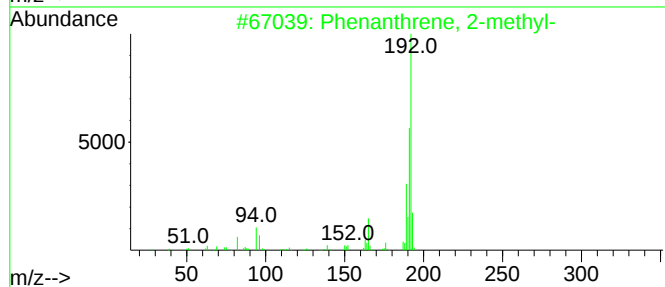
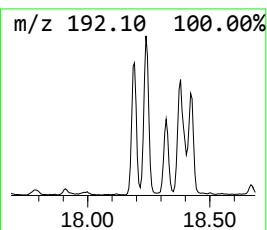
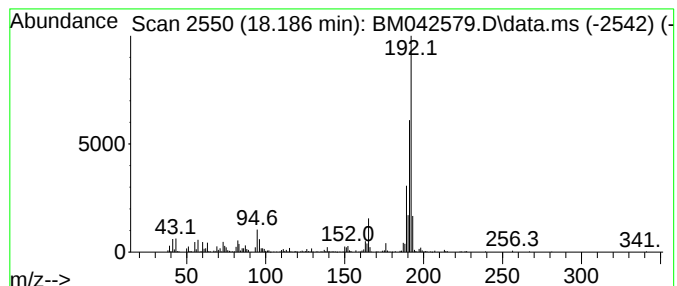
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Phenanthrene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.186	16.28 ng	536851	Phenanthrene-d10	17.321

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
3		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	94
4		Anthracene, 2-methyl-	192	C15H12	000613-12-7	94
5		Anthracene, 1-methyl-	192	C15H12	000610-48-0	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110323\
Data File : BM042579.D
Acq On : 03 Nov 2023 20:23
Operator : MA/JU
Sample : 05220-02 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
Z-5(0-5)

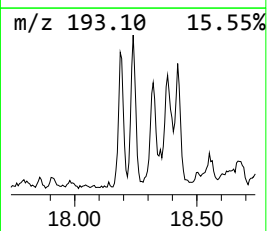
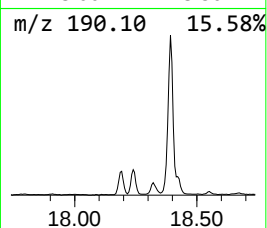
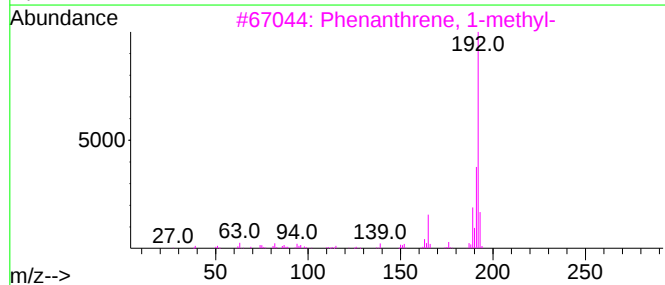
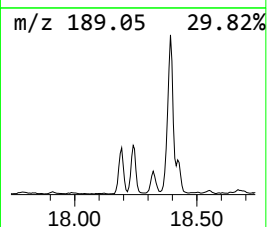
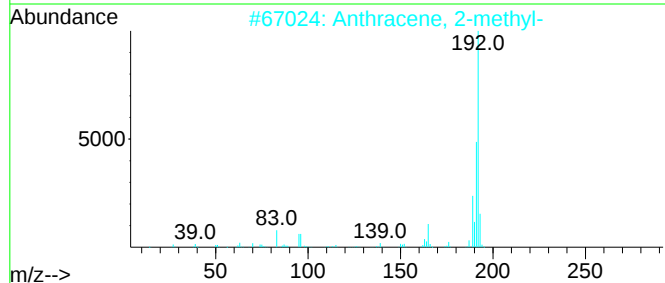
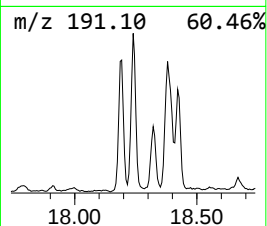
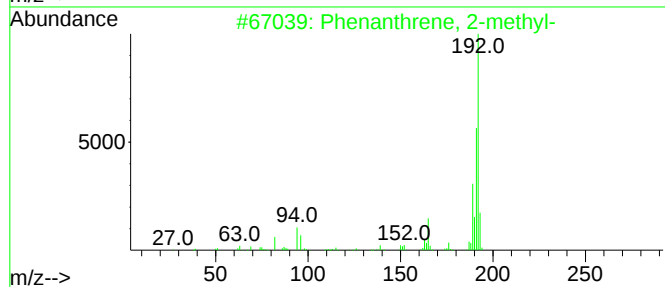
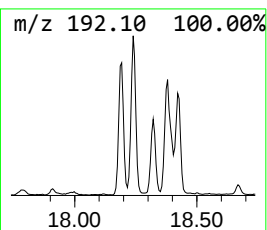
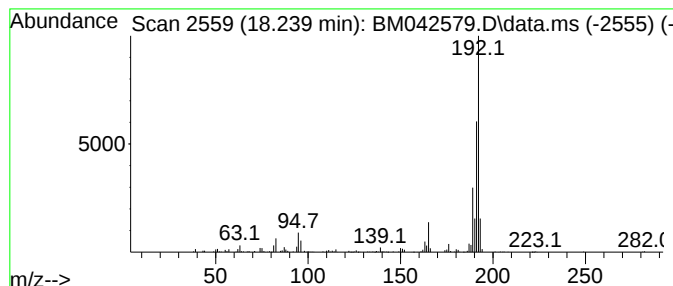
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Anthracene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.239	11.44 ng	377284	Phenanthrene-d10	17.321

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
2			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
3			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	94
4			Anthracene, 9-methyl-	192	C15H12	000779-02-2	93
5			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110323\
Data File : BM042579.D
Acq On : 03 Nov 2023 20:23
Operator : MA/JU
Sample : 05220-02 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
Z-5(0-5)

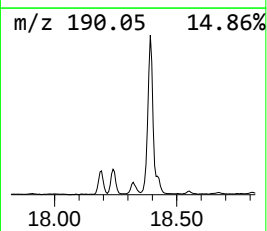
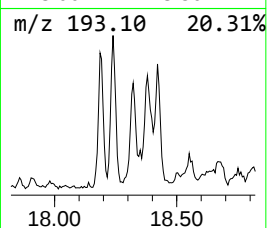
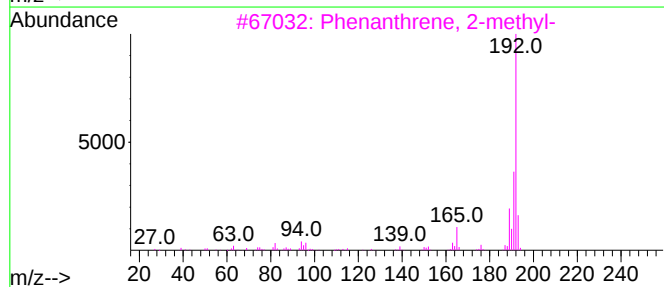
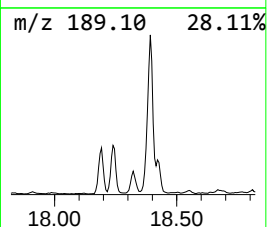
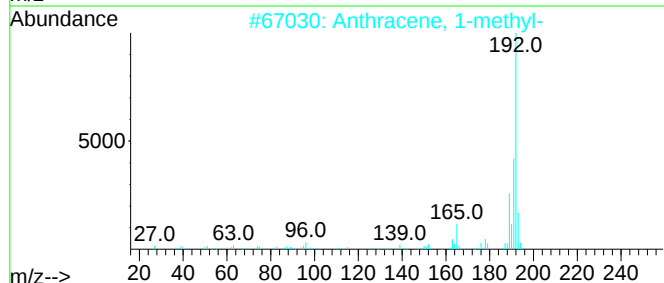
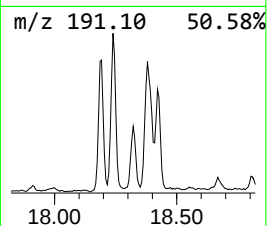
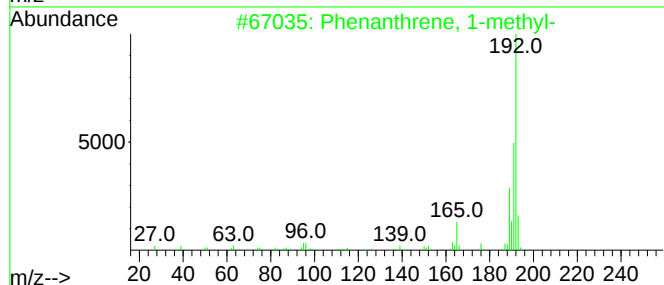
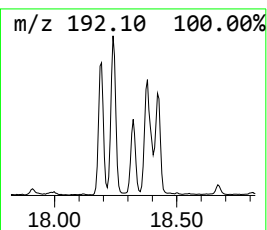
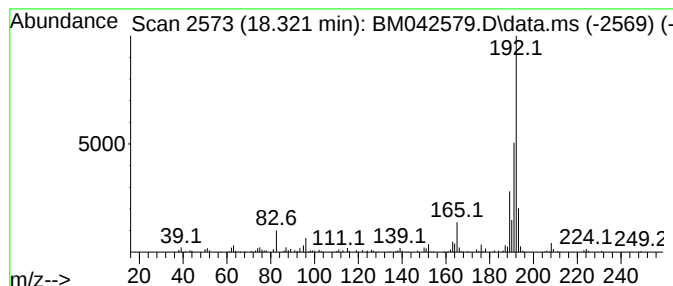
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Anthracene, 1-methyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.321	4.88 ng	160907	Phenanthrene-d10	17.321

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
2		Anthracene, 1-methyl-	192	C15H12	000610-48-0	96
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
4		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	94
5		Anthracene, 9-methyl-	192	C15H12	000779-02-2	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110323\
Data File : BM042579.D
Acq On : 03 Nov 2023 20:23
Operator : MA/JU
Sample : 05220-02 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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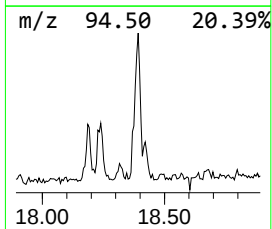
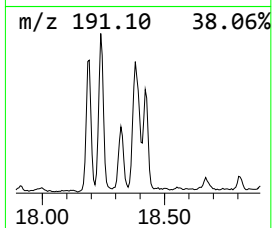
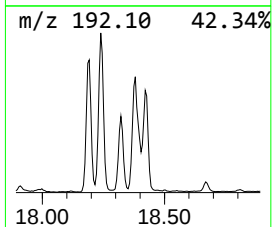
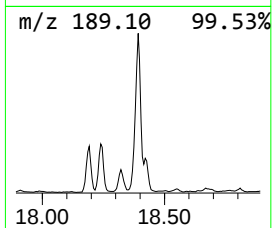
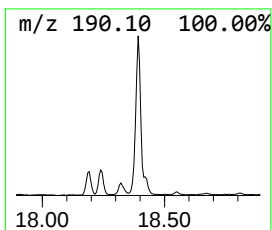
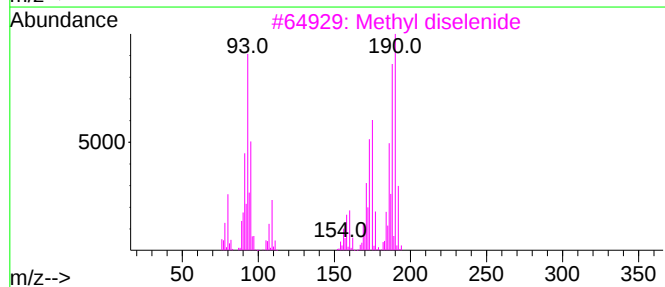
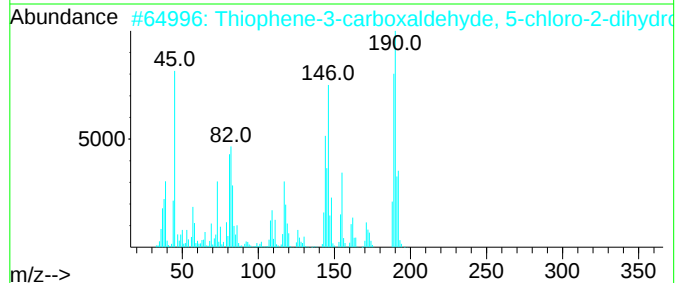
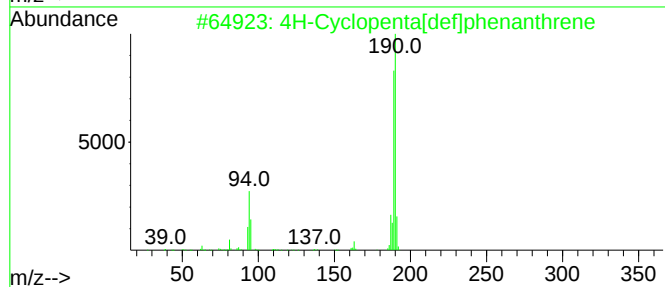
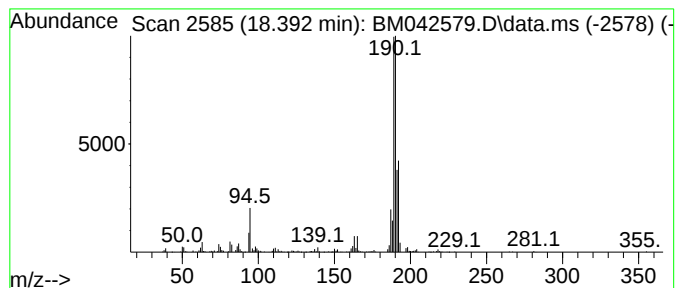
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 4H-Cyclopenta[def]phenanthrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.392	19.29 ng	636265	Phenanthrene-d10	17.321

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	76
2		Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	53
3		Methyl diselenide	190	C2H6Se2	007101-31-7	45
4		6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	42
5		2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110323\
Data File : BM042579.D
Acq On : 03 Nov 2023 20:23
Operator : MA/JU
Sample : 05220-02 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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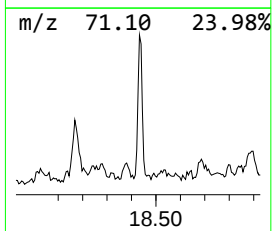
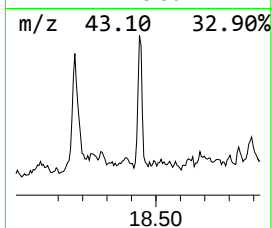
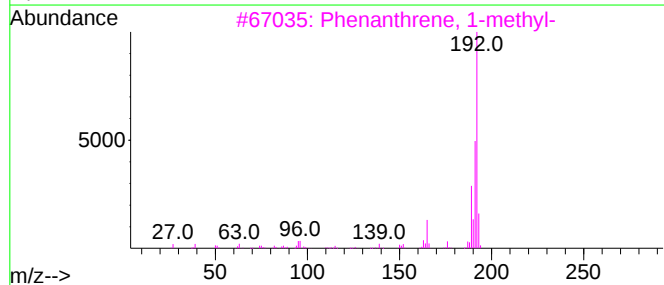
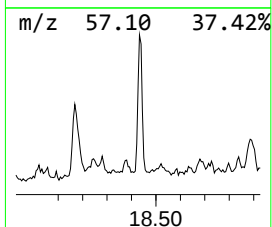
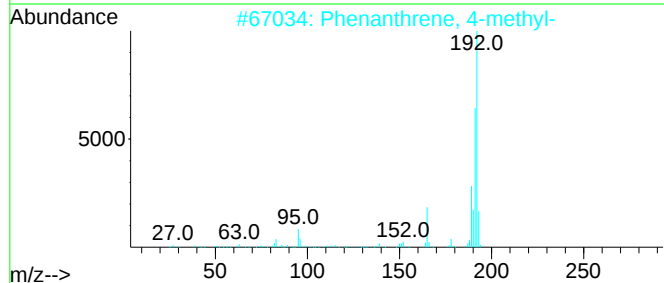
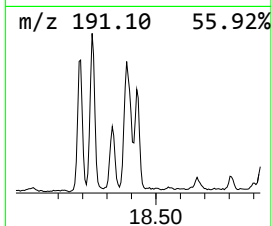
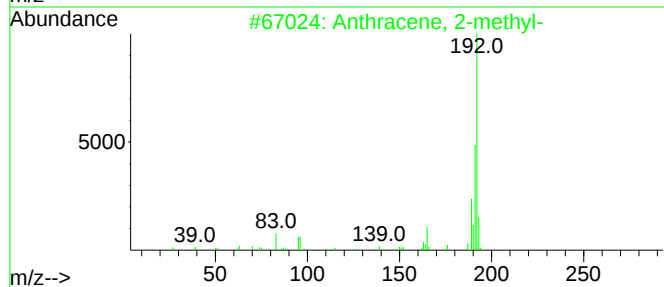
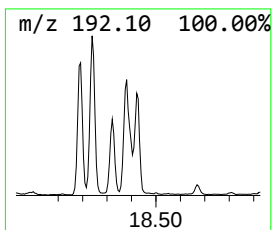
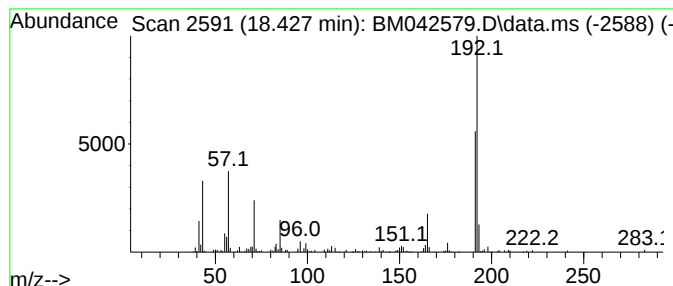
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Phenanthrene, 4-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.427	8.30 ng	273672	Phenanthrene-d10	17.321

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 2-methyl-	192	C15H12	000613-12-7	76
2		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	70
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	70
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	68
5		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110323\
Data File : BM042579.D
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Sample : 05220-02 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

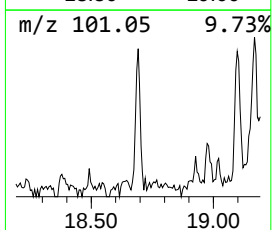
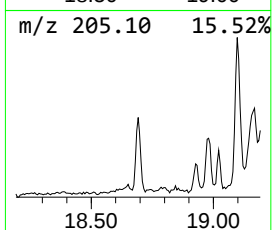
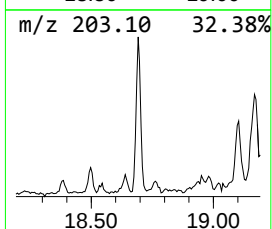
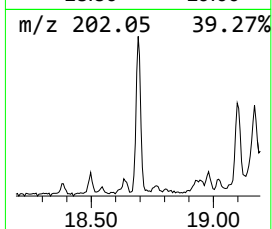
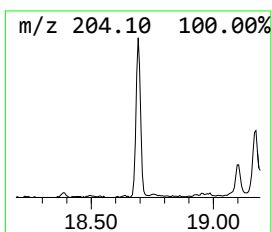
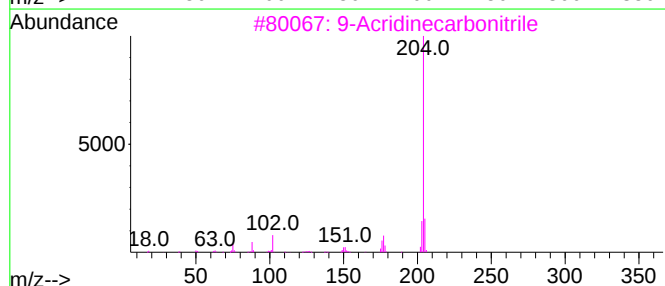
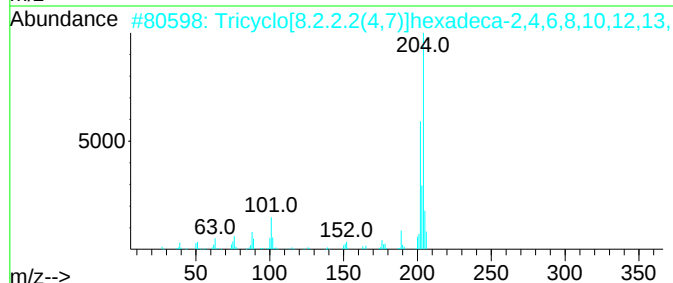
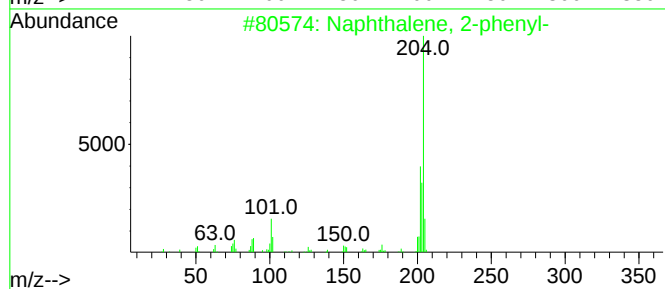
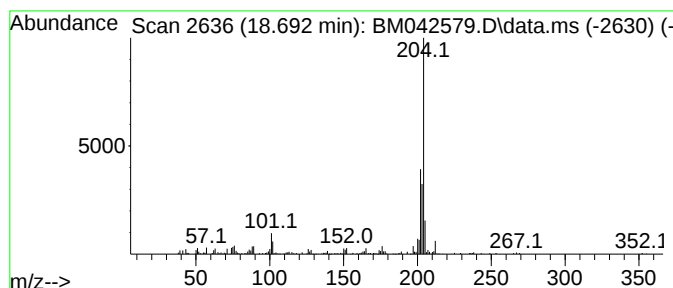
Instrument :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Naphthalene, 2-phenyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.692	6.66 ng	219631	Phenanthrene-d10	17.321	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	86
2	Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	62
3	9-Acridinecarbonitrile	204	C14H8N2	005326-19-2	55
4	[1,1'-Biphenyl]-2-ol, 5-chloro-	204	C12H9ClO	000607-12-5	49
5	9,10-ethenoanthracene, 9,10-dihy...	204	C16H12	1000400-47-8	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110323\
Data File : BM042579.D
Acq On : 03 Nov 2023 20:23
Operator : MA/JU
Sample : 05220-02 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

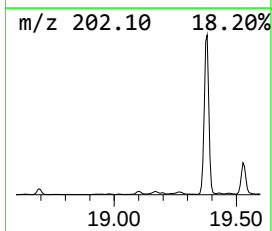
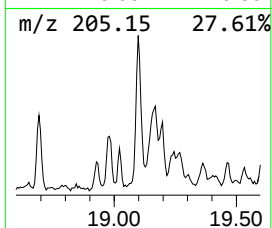
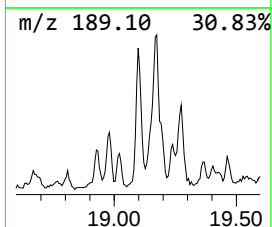
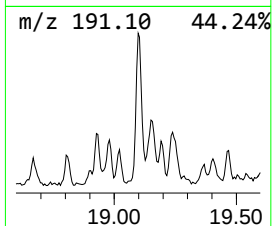
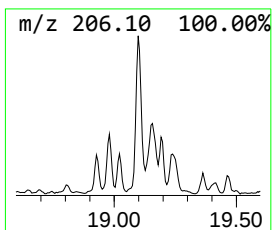
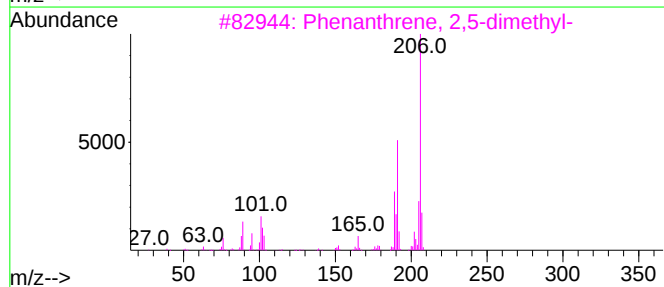
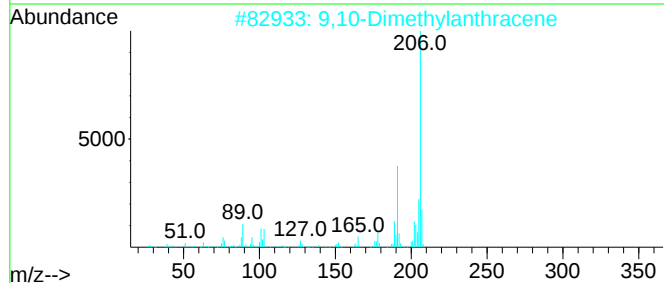
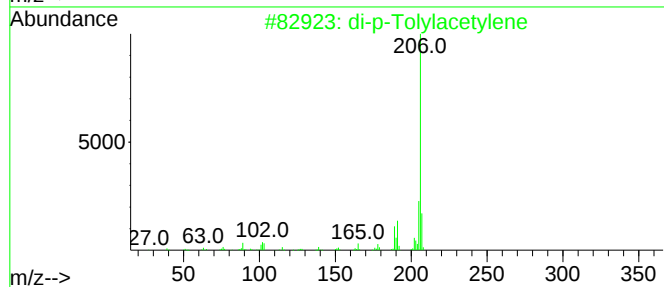
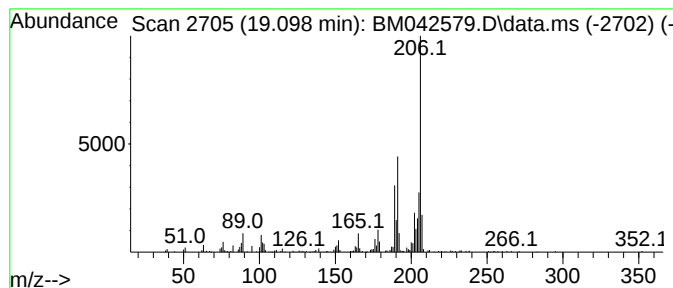
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 di-p-Tolylacetylene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.098	6.24 ng	205686	Phenanthrene-d10	17.321	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	di-p-Tolylacetylene	206	C16H14	002789-88-0	94
2	9,10-Dimethylantracene	206	C16H14	000781-43-1	91
3	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	91
4	Naphthalene, 1,2-dihydro-4-phenyl-	206	C16H14	007469-40-1	90
5	Naphthalene, 1,2-dihydro-1-phenyl-	206	C16H14	016606-46-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110323\
Data File : BM042579.D
Acq On : 03 Nov 2023 20:23
Operator : MA/JU
Sample : 05220-02 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
Z-5(0-5)

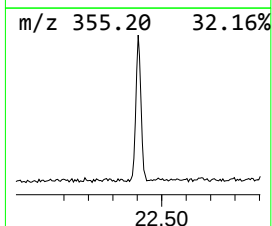
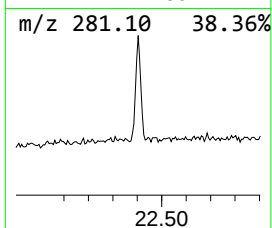
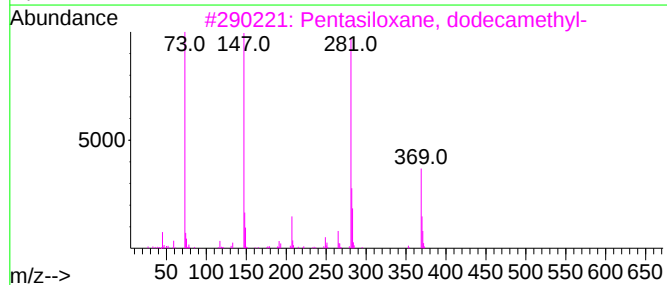
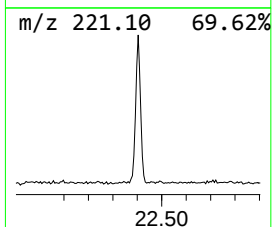
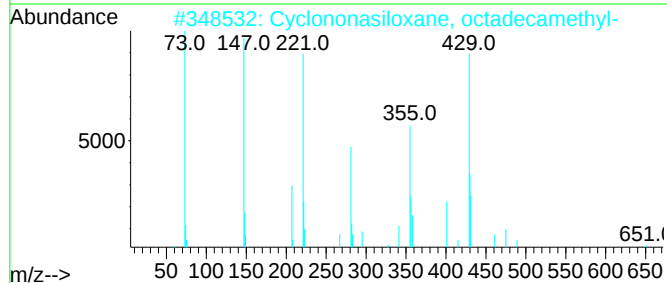
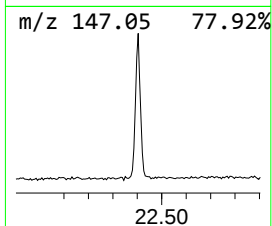
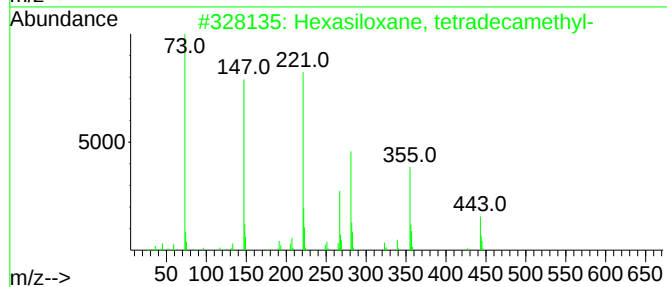
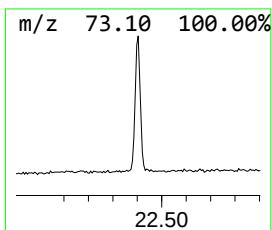
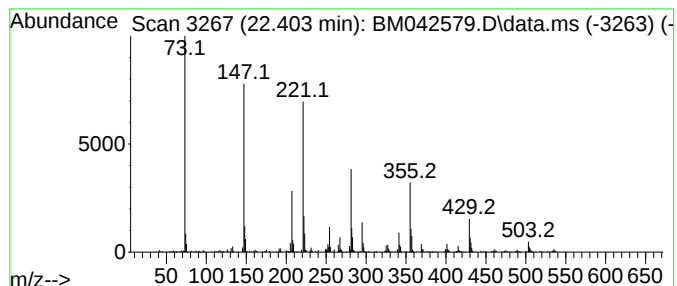
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 Hexasiloxane, tetradecamethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.403	7.57 ng	666707	Chrysene-d12	21.503

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexasiloxane, tetradecamethyl-	458	C14H42O5Si6	000107-52-8	83
2			Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	78
3			Pentasiloxane, dodecamethyl-	384	C12H36O4Si5	000141-63-9	43
4			N-Benzyl-N-ethyl-p-isopropylbenz...	281	C19H23NO	015089-22-2	43
5			Mercaptoacetic acid, 2TMS deriva...	236	C8H20O2SSi2	006398-62-5	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110323\
Data File : BM042579.D
Acq On : 03 Nov 2023 20:23
Operator : MA/JU
Sample : 05220-02 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
Z-5(0-5)

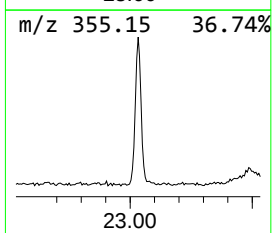
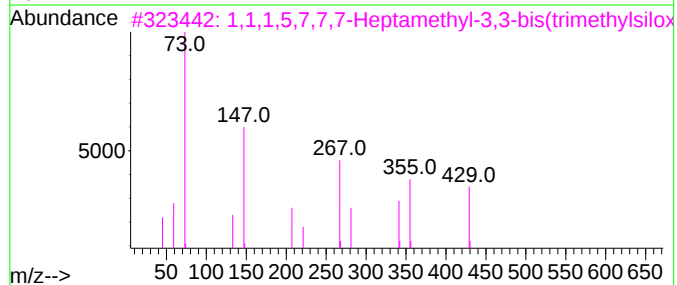
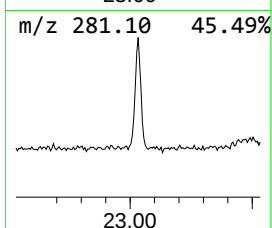
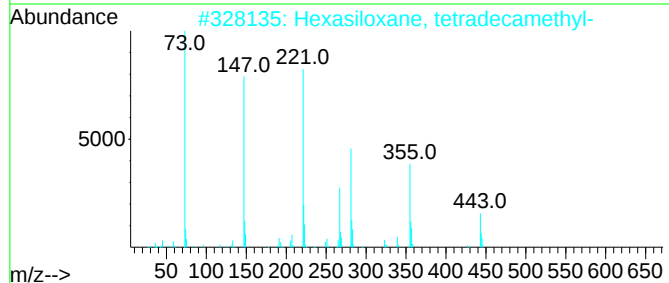
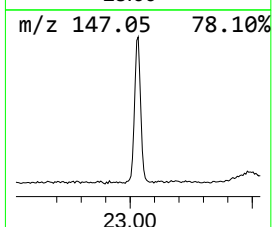
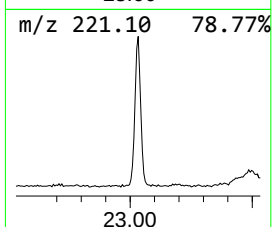
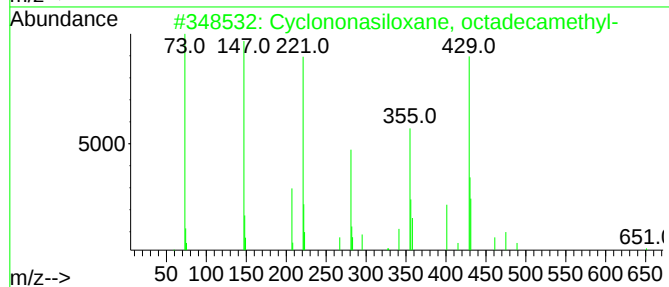
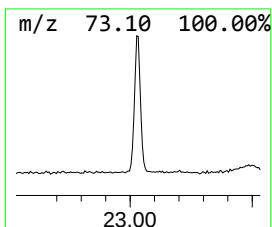
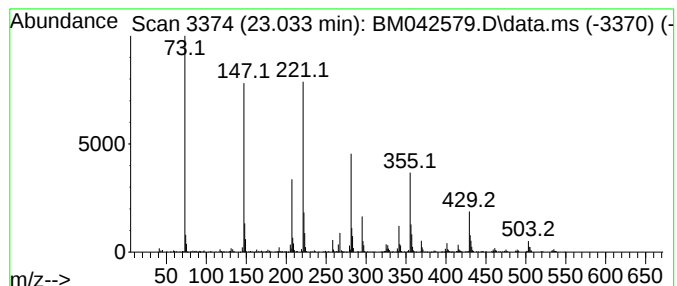
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 Cyclononasiloxane, octadeca... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.033	21.67 ng	619020	Perylene-d12	23.927

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	91
2			Hexasiloxane, tetradecamethyl-	458	C14H42O5Si6	000107-52-8	80
3			1,1,1,5,7,7,7-Heptamethyl-3,3-bi...	444	C13H40O5Si6	038147-00-1	43
4			3,6-Dioxa-2,4,5,7-tetrasilaoctan...	294	C10H30O2Si4	004342-25-0	40
5			Mercaptoacetic acid, 2TMS deriva...	236	C8H20O2SSi2	006398-62-5	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110323\
Data File : BM042579.D
Acq On : 03 Nov 2023 20:23
Operator : MA/JU
Sample : 05220-02 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
Z-5(0-5)

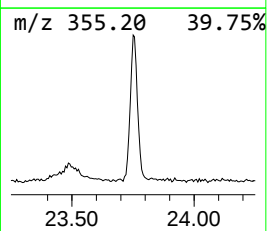
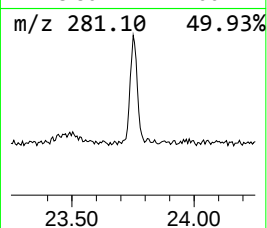
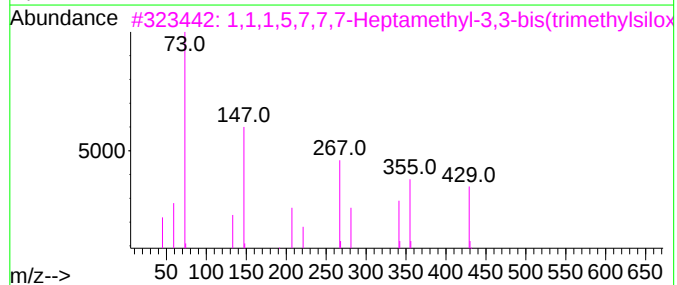
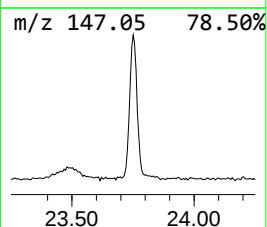
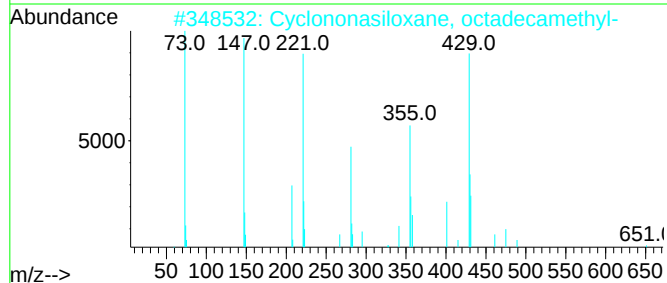
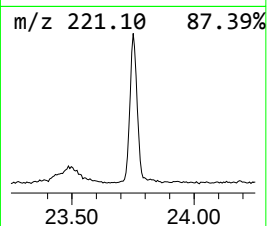
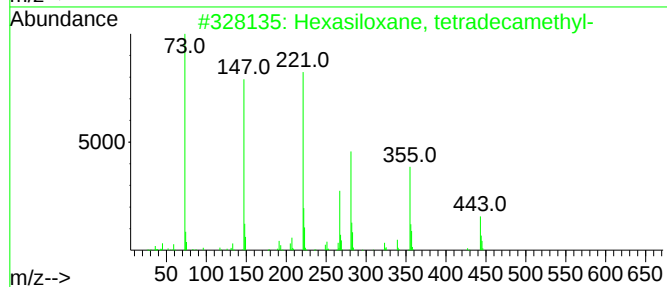
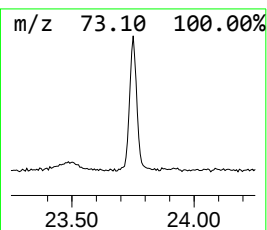
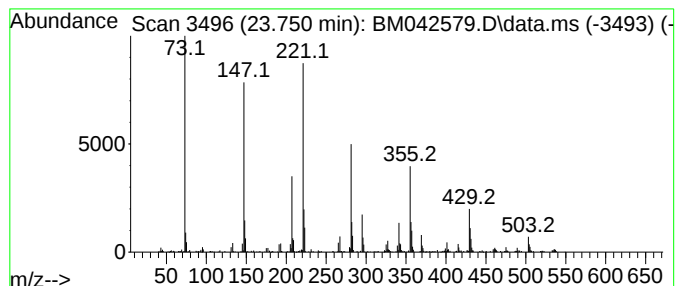
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 unknown23.750 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.750	22.49 ng	642419	Perylene-d12	23.927

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexasiloxane, tetradecamethyl-	458	C14H42O5Si6	000107-52-8	83
2			Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	72
3			1,1,1,5,7,7,7-Heptamethyl-3,3-bi...	444	C13H40O5Si6	038147-00-1	47
4			3,6-Dioxa-2,4,5,7-tetrasilaoctan...	294	C10H30O2Si4	004342-25-0	38
5			Mercaptoacetic acid, 2TMS deriva...	236	C8H20O2SSi2	006398-62-5	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110323\
Data File : BM042579.D
Acq On : 03 Nov 2023 20:23
Operator : MA/JU
Sample : 05220-02 5X
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
Z-5(0-5)

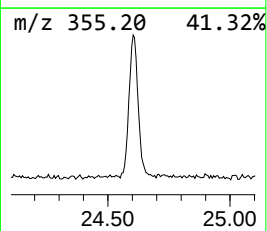
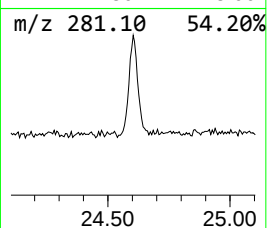
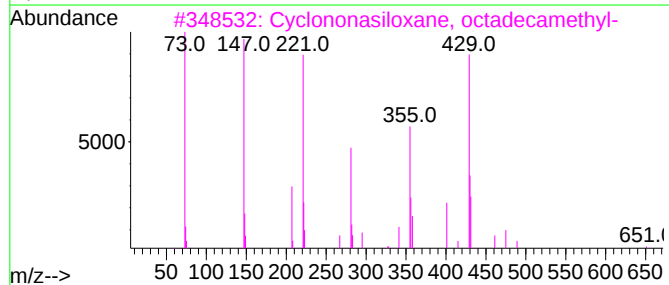
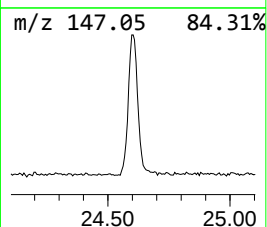
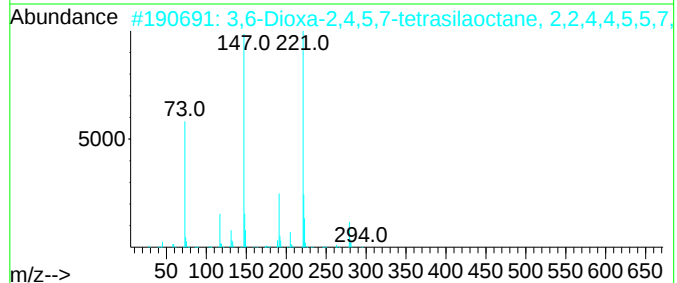
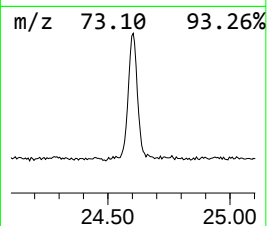
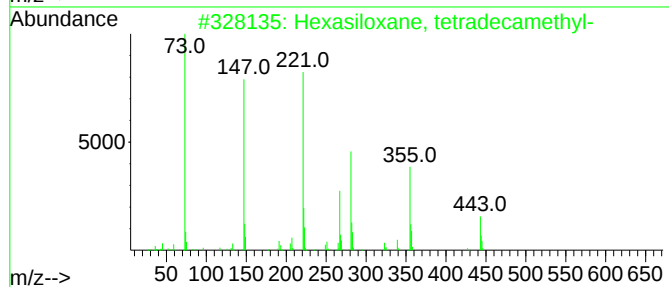
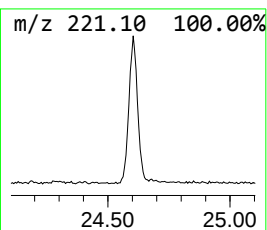
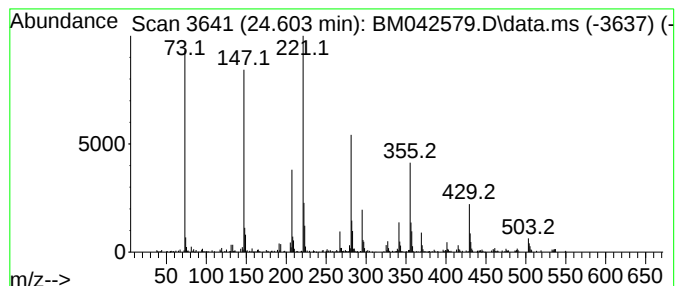
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 unknown24.603 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.603	28.33 ng	809194	Perylene-d12	23.927

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexasiloxane, tetradecamethyl-	458	C14H42O5Si6	000107-52-8	76
2			3,6-Dioxa-2,4,5,7-tetrasilaoctan...	294	C10H30O2Si4	004342-25-0	64
3			Cyclononasiloxane, octadecamethyl-	666	C18H54O9Si9	000556-71-8	64
4			Pentasiloxane, dodecamethyl-	384	C12H36O4Si5	000141-63-9	25
5			Phthalic acid, 2TMS derivative	310	C14H22O4Si2	002078-22-0	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110323\
 Data File : BM042579.D
 Acq On : 03 Nov 2023 20:23
 Operator : MA/JU
 Sample : 05220-02 5X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM103023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Naphthalene, 2,...	13.722	7.7	ng	213588	3	14.569	552299	20.0
Naphthalene, 2,...	13.880	9.8	ng	270618	3	14.569	552299	20.0
Naphthalene, 1,...	14.116	4.3	ng	117363	3	14.569	552299	20.0
Naphthalene, 2,...	14.945	5.1	ng	140750	3	14.569	552299	20.0
Naphthalene, 1,...	15.163	4.4	ng	121909	3	14.569	552299	20.0
9H-Fluorene, 2-...	16.610	4.6	ng	151691	4	17.321	659558	20.0
Dibenzothiophene	17.139	6.6	ng	216839	4	17.321	659558	20.0
4-Methylnaphtho...	17.898	5.7	ng	189318	4	17.321	659558	20.0
1-Methyldibenzo...	18.045	4.8	ng	156968	4	17.321	659558	20.0
Phenanthrene, 2...	18.186	16.3	ng	536851	4	17.321	659558	20.0
Anthracene, 2-m...	18.239	11.4	ng	377284	4	17.321	659558	20.0
Anthracene, 1-m...	18.321	4.9	ng	160907	4	17.321	659558	20.0
4H-Cyclopenta[d...	18.392	19.3	ng	636265	4	17.321	659558	20.0
Phenanthrene, 4...	18.427	8.3	ng	273672	4	17.321	659558	20.0
Naphthalene, 2-...	18.692	6.7	ng	219631	4	17.321	659558	20.0
di-p-Tolylacety...	19.098	6.2	ng	205686	4	17.321	659558	20.0
Hexasiloxane, t...	22.403	7.6	ng	666707	5	21.503	1762240	20.0
Cyclononasiloxa...	23.033	21.7	ng	619020	6	23.927	571321	20.0
unknown23.750	23.750	22.5	ng	642419	6	23.927	571321	20.0
unknown24.603	24.603	28.3	ng	809194	6	23.927	571321	20.0