

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021422\
 Data File : BP009148.D
 Acq On : 14 Feb 2022 22:47
 Operator : CG/JU
 Sample : N1526-03 2X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 HE-2-021122

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-BP020722.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP009148.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.999	168	172	182	rVB	32721	57636	1.11%	0.099%
2	5.393	402	409	434	rBV	1665756	2894866	55.87%	4.983%
3	5.811	473	480	488	rVB5	38741	73719	1.42%	0.127%
4	6.064	515	523	531	rBV4	24460	62165	1.20%	0.107%
5	6.928	665	670	675	rBV5	33003	77746	1.50%	0.134%
6	6.993	675	681	697	rVV	1310331	2660624	51.35%	4.580%
7	7.111	697	701	707	rVB6	25353	54359	1.05%	0.094%
8	7.275	725	729	737	rVB2	27791	53263	1.03%	0.092%
9	7.416	748	753	765	rBV3	62134	129451	2.50%	0.223%
10	7.793	807	817	825	rBV	930693	1640453	31.66%	2.824%
11	8.075	857	865	872	rVB5	34288	77274	1.49%	0.133%
12	8.340	897	910	915	rBV8	52453	154894	2.99%	0.267%
13	8.899	994	1005	1008	rBV8	34916	84257	1.63%	0.145%
14	8.958	1008	1015	1023	rVV	879087	1734131	33.47%	2.985%
15	9.016	1023	1025	1036	rVB	85643	153258	2.96%	0.264%
16	9.146	1042	1047	1054	rVB8	27405	61995	1.20%	0.107%
17	9.499	1100	1107	1114	rVB5	21180	55836	1.08%	0.096%
18	10.593	1282	1293	1298	rBV	1106469	2075207	40.05%	3.572%
19	10.640	1298	1301	1309	rVB	177003	317665	6.13%	0.547%
20	12.257	1568	1576	1585	rBV	117314	213783	4.13%	0.368%
21	12.475	1604	1613	1623	rBV	133984	264167	5.10%	0.455%
22	13.057	1704	1712	1728	rBV	2554728	4127976	79.68%	7.106%
23	13.269	1743	1748	1754	rVB2	44158	75498	1.46%	0.130%
24	13.610	1799	1806	1813	rBV3	90997	245928	4.75%	0.423%
25	13.757	1825	1831	1837	rBV	156977	308300	5.95%	0.531%
26	13.816	1837	1841	1845	rVV	78566	121699	2.35%	0.209%
27	13.898	1851	1855	1863	rBV2	42940	89726	1.73%	0.154%
28	13.998	1867	1872	1875	rBV	60012	93856	1.81%	0.162%
29	14.163	1894	1900	1912	rBV	224239	391734	7.56%	0.674%
30	14.340	1923	1930	1936	rBV3	47856	93717	1.81%	0.161%
31	14.434	1938	1946	1953	rVV	1505319	2503680	48.32%	4.310%
32	14.498	1953	1957	1966	rVB	155399	293568	5.67%	0.505%
33	14.657	1978	1984	1992	rVB4	46530	112550	2.17%	0.194%
34	14.840	2010	2015	2021	rVB	130863	212867	4.11%	0.366%
35	14.916	2024	2028	2033	rVB	70009	102757	1.98%	0.177%
36	15.057	2045	2052	2057	rBV3	81606	188175	3.63%	0.324%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M

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37	15.110	2058	2061	2065	rVB	43590	62024	1.20%	0.107%
38	15.228	2074	2081	2085	rBV5	56979	135463	2.61%	0.233%
39	15.493	2120	2126	2137	rVB	253678	493239	9.52%	0.849%
40	15.616	2138	2147	2149	rBV2	46584	95256	1.84%	0.164%
41	15.698	2157	2161	2166	rBV2	52669	79101	1.53%	0.136%
42	15.787	2171	2176	2182	rVV3	49927	92798	1.79%	0.160%
43	15.940	2193	2202	2222	rBV	1813658	3142229	60.65%	5.409%
44	16.175	2235	2242	2247	rBV5	88481	189320	3.65%	0.326%
45	16.498	2290	2297	2302	rBV2	104738	236011	4.56%	0.406%
46	16.551	2302	2306	2312	rVB2	72717	121942	2.35%	0.210%
47	16.645	2319	2322	2328	rVB2	58578	88774	1.71%	0.153%
48	16.857	2355	2358	2364	rVB3	54211	86918	1.68%	0.150%
49	17.016	2380	2385	2393	rVB	184519	327167	6.31%	0.563%
50	17.192	2410	2415	2419	rBV	1878442	2877659	55.54%	4.954%
51	17.239	2419	2423	2429	rVB	1736842	2389582	46.12%	4.113%
52	17.334	2434	2439	2444	rVB	482520	715366	13.81%	1.231%
53	17.616	2482	2487	2496	rBV	137933	297091	5.73%	0.511%
54	17.781	2511	2515	2523	rVB2	140560	226263	4.37%	0.389%
55	17.869	2526	2530	2534	rVV	199428	253365	4.89%	0.436%
56	17.922	2535	2539	2545	rVB	91733	162443	3.14%	0.280%
57	18.069	2559	2564	2567	rBV	386430	608263	11.74%	1.047%
58	18.116	2568	2572	2579	rVB	539301	786189	15.17%	1.353%
59	18.192	2581	2585	2589	rVV	196984	292117	5.64%	0.503%
60	18.257	2590	2596	2599	rVV2	550448	1010220	19.50%	1.739%
61	18.292	2599	2602	2607	rVB	298399	416183	8.03%	0.716%
62	18.551	2642	2646	2650	rBV2	184902	257550	4.97%	0.443%
63	18.763	2679	2682	2684	rBV2	95896	118877	2.29%	0.205%
64	18.839	2692	2695	2698	rBV	146548	158198	3.05%	0.272%
65	18.969	2711	2717	2720	rBV2	609957	1135580	21.92%	1.955%
66	19.004	2720	2723	2728	rVB4	472460	693104	13.38%	1.193%
67	19.210	2754	2758	2764	rBV	1892831	2594721	50.08%	4.467%
68	19.563	2810	2818	2827	rBV	1966138	3524424	68.03%	6.067%
69	19.757	2847	2851	2856	rVB	4059057	5181002	100.00%	8.919%
70	20.145	2915	2917	2920	rVB	270859	269473	5.20%	0.464%
71	21.292	3107	3112	3116	rBV2	2645619	4189148	80.86%	7.211%
72	23.698	3516	3521	3527	rVB2	1632705	2925378	56.46%	5.036%

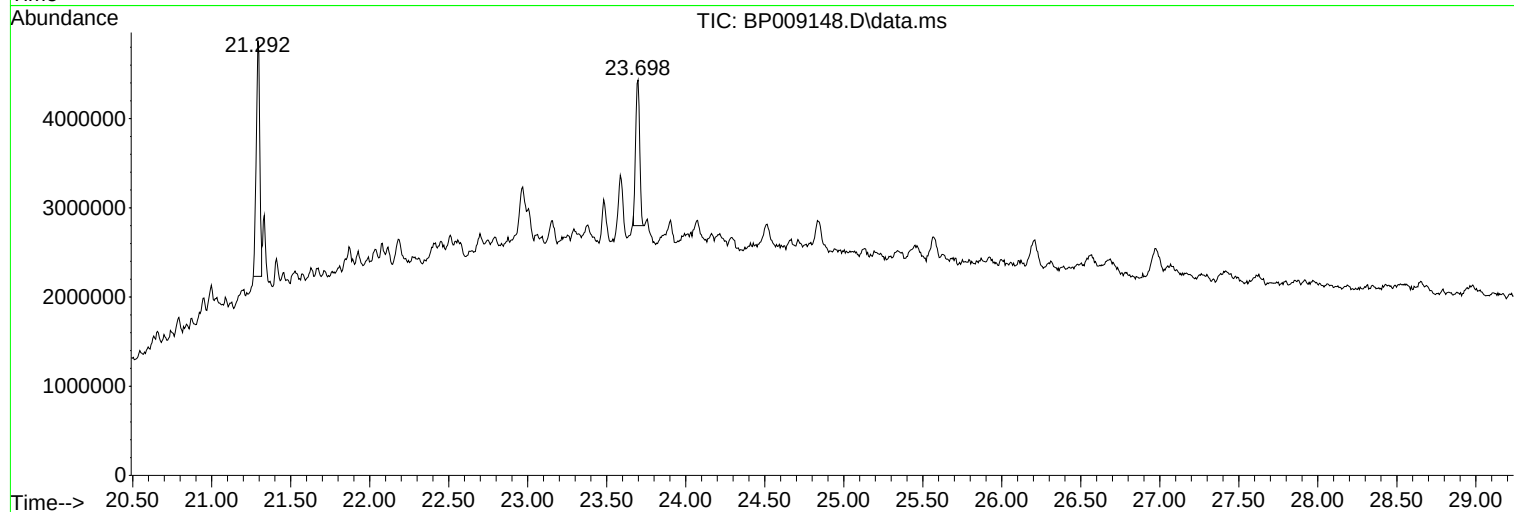
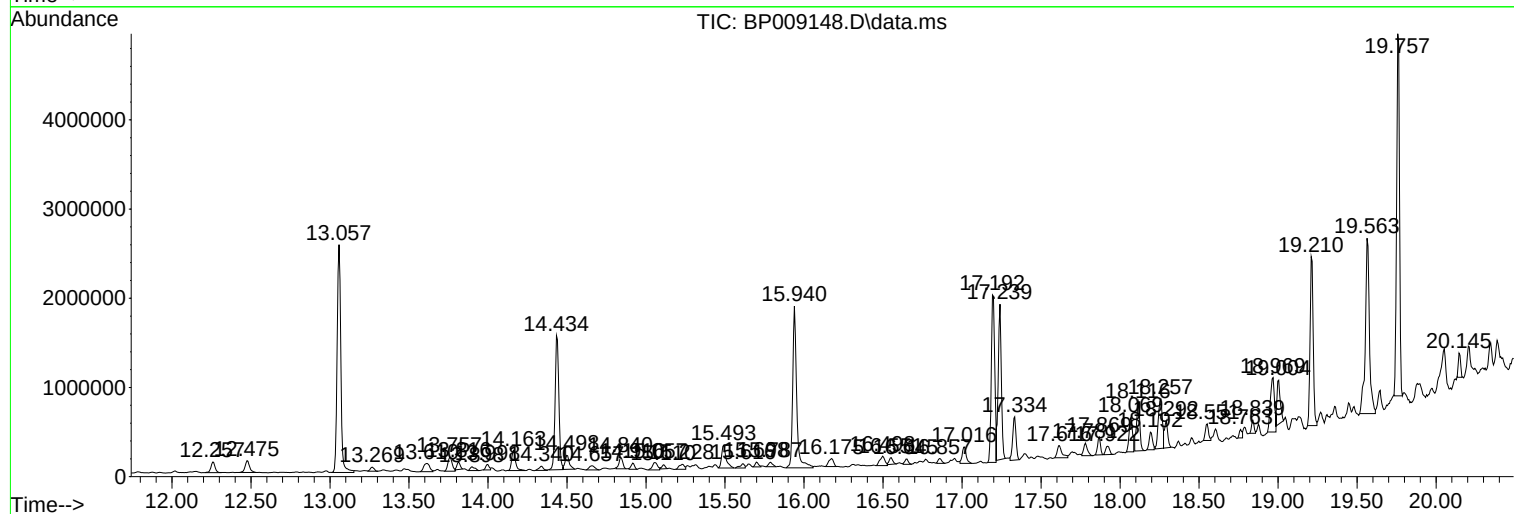
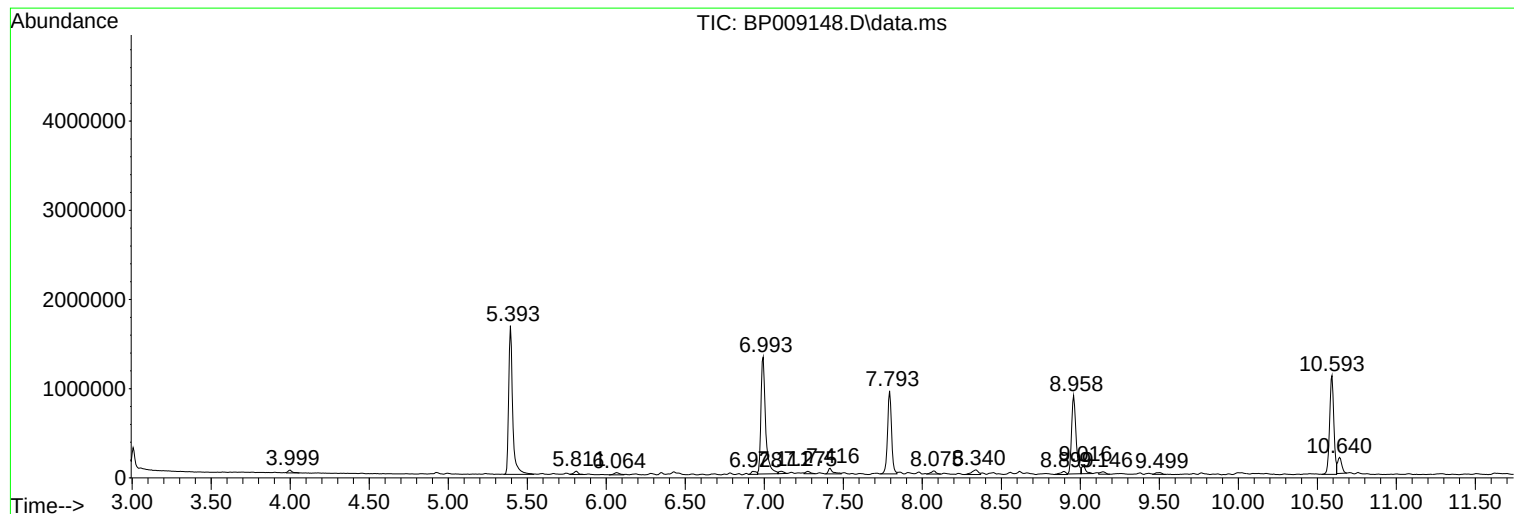
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Sample : N1526-03 2X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
HE-2-021122

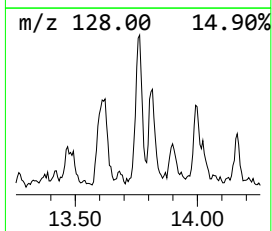
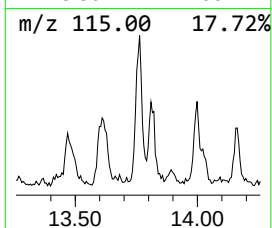
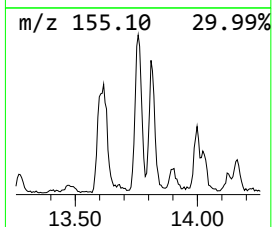
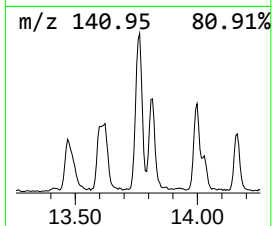
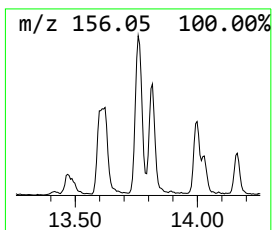
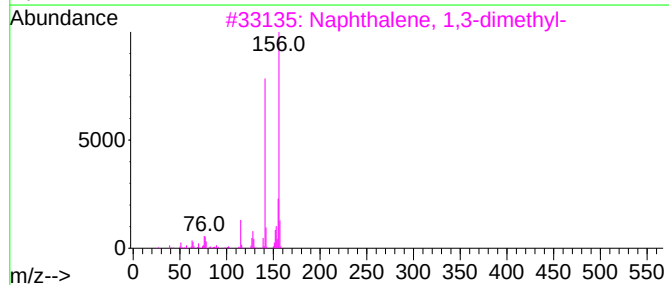
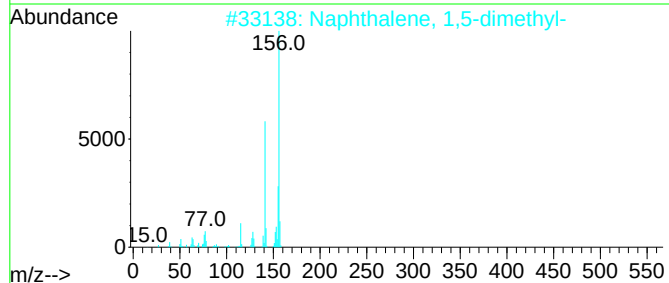
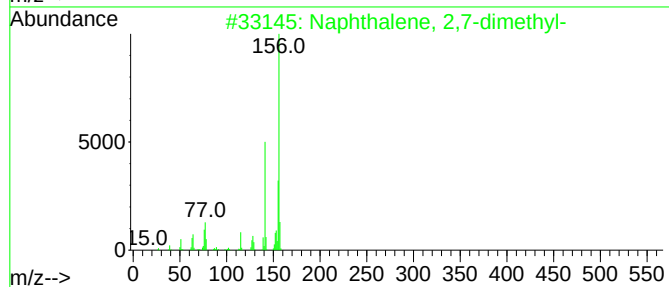
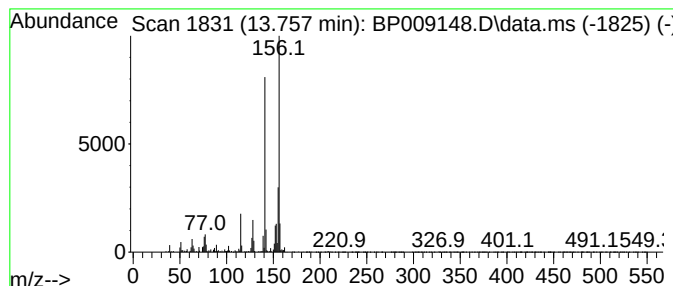
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.757	2.46 ng	308300	Acenaphthene-d10	14.434

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
2			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
3			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
4			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
5			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96



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Instrument :
BNA_P
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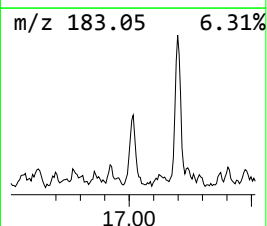
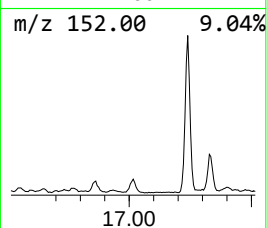
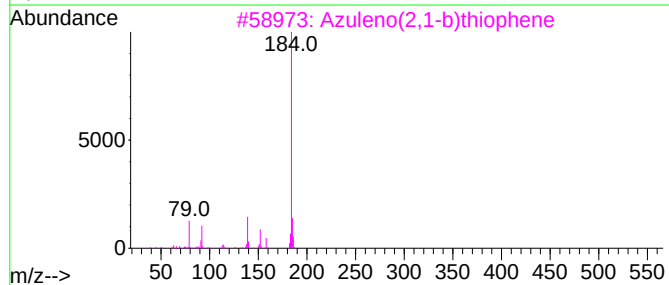
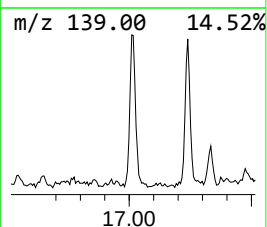
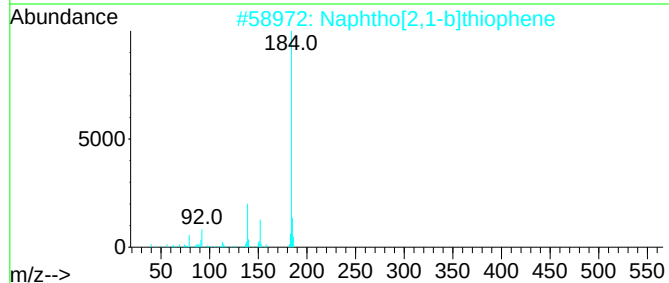
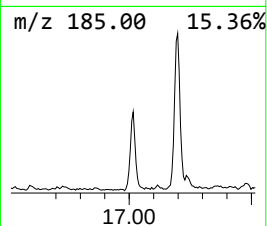
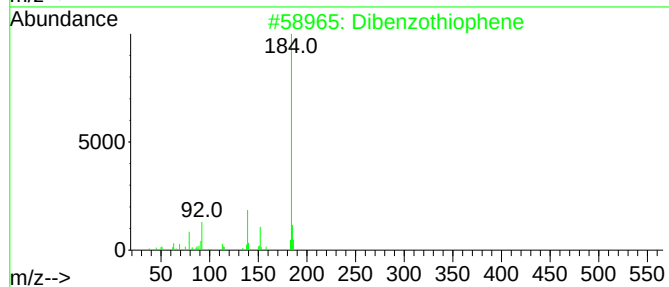
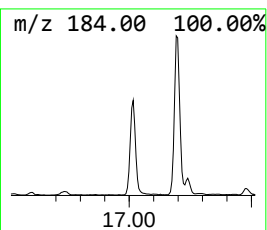
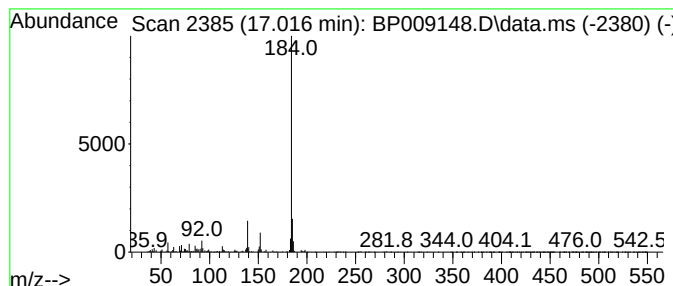
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Dibenzothiophene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.016	2.27 ng	327167	Phenanthrene-d10	17.192

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzothiophene	184	C12H8S	000132-65-0	95
2			Naphtho[2,1-b]thiophene	184	C12H8S	000233-02-3	94
3			Azuleno(2,1-b)thiophene	184	C12H8S	000248-13-5	91
4			Naphtho[2,3-b]thiophene	184	C12H8S	000268-77-9	90
5			Thianthrene, 5-oxide	232	C12H8OS2	002362-50-7	78



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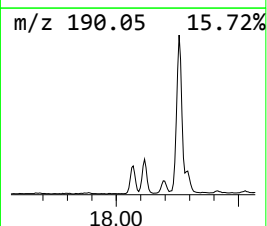
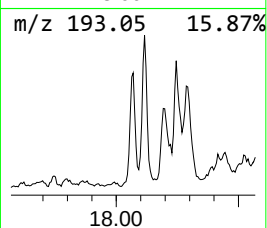
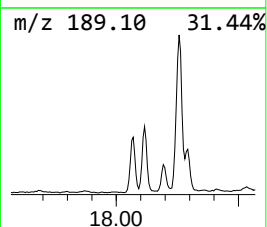
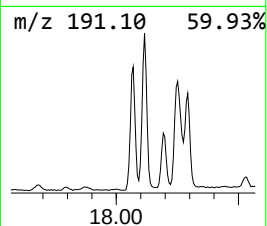
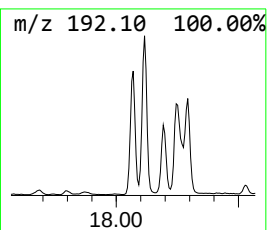
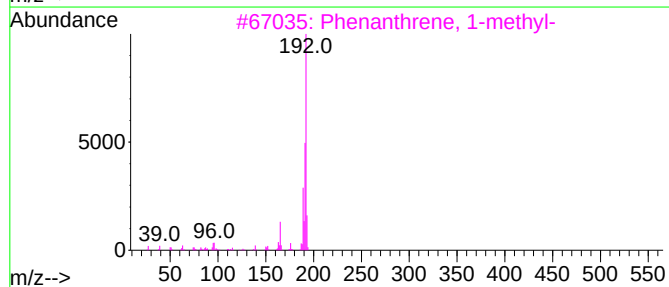
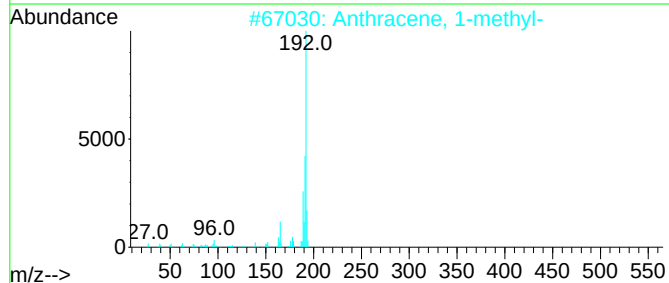
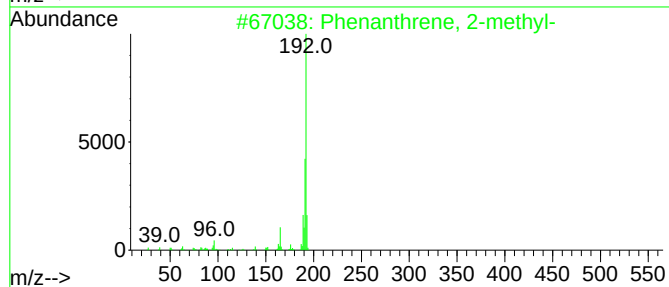
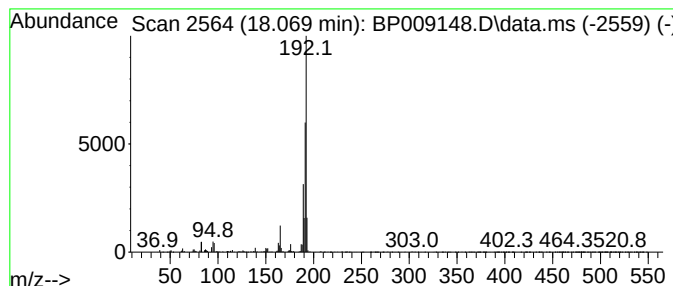
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Phenanthrene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.069	4.23 ng	608263	Phenanthrene-d10	17.192

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	93
2		Anthracene, 1-methyl-	192	C15H12	000610-48-0	93
3		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93
4		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	93
5		Anthracene, 2-methyl-	192	C15H12	000613-12-7	93



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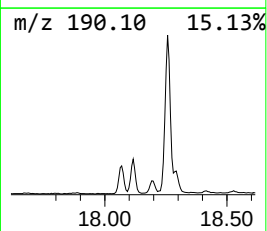
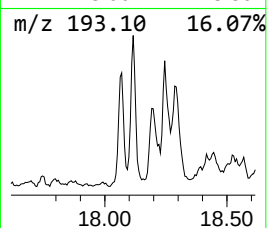
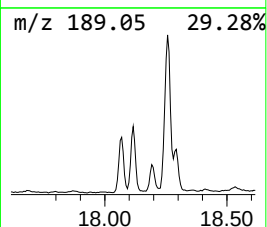
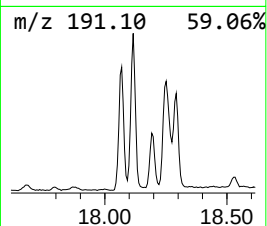
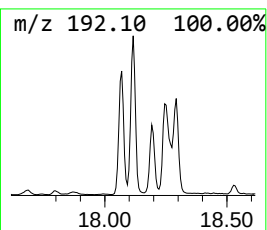
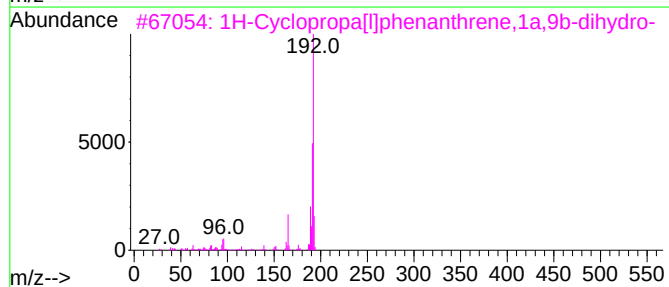
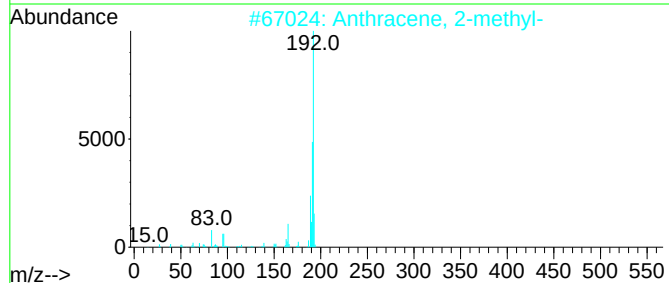
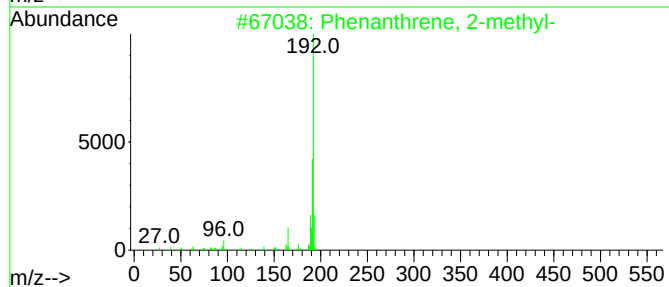
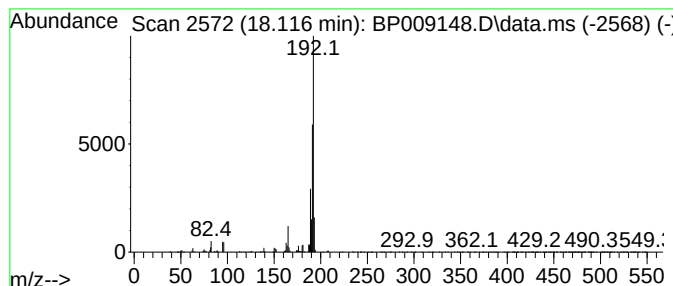
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Anthracene, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.116	5.46 ng	786189	Phenanthrene-d10	17.192

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		1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	96
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	95
5		Anthracene, 1-methyl-	192	C15H12	000610-48-0	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021422\
Data File : BP009148.D
Acq On : 14 Feb 2022 22:47
Operator : CG/JU
Sample : N1526-03 2X
Misc :
ALS Vial : 19 Sample Multiplier: 1

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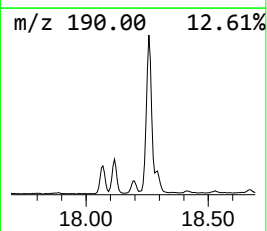
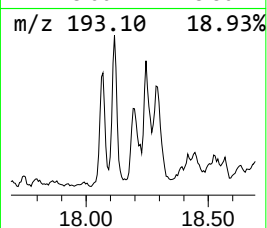
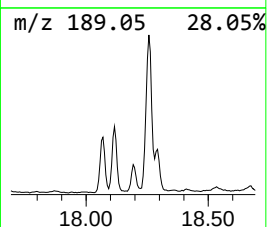
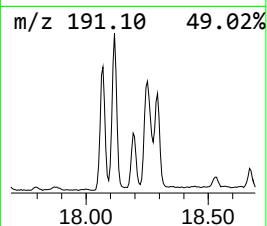
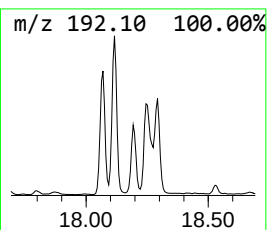
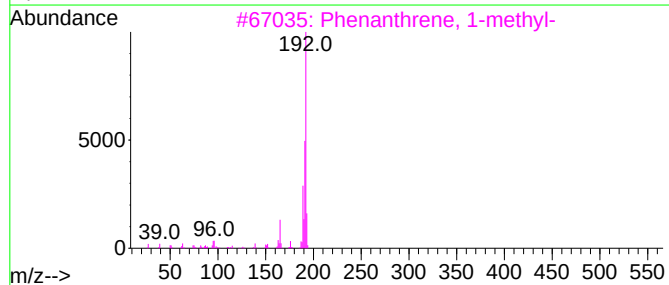
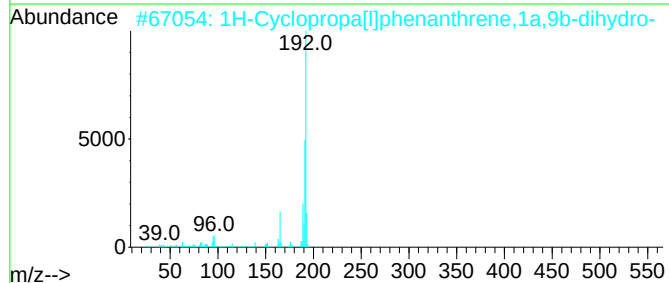
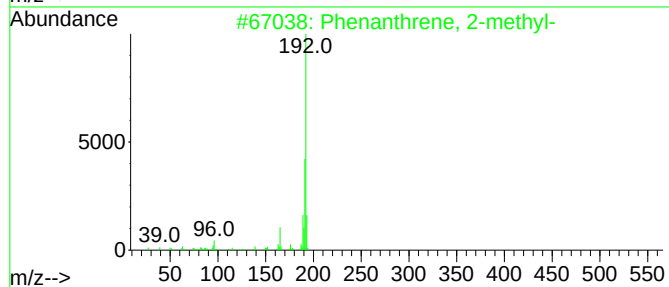
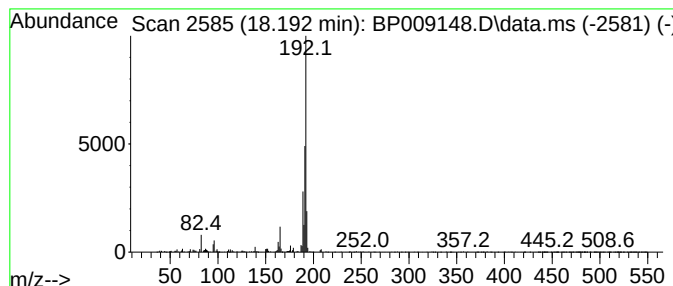
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 1H-Cyclopropa[l]phenanthren... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.192	2.03 ng	292117	Phenanthrene-d10	17.192

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021422\
Data File : BP009148.D
Acq On : 14 Feb 2022 22:47
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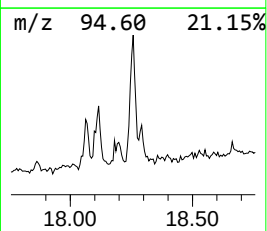
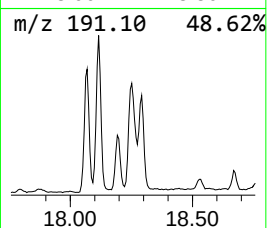
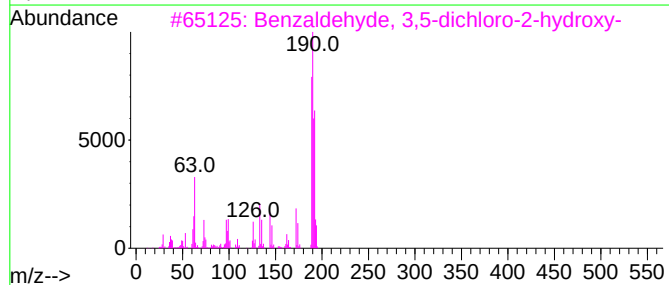
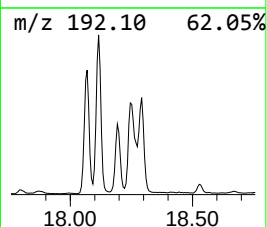
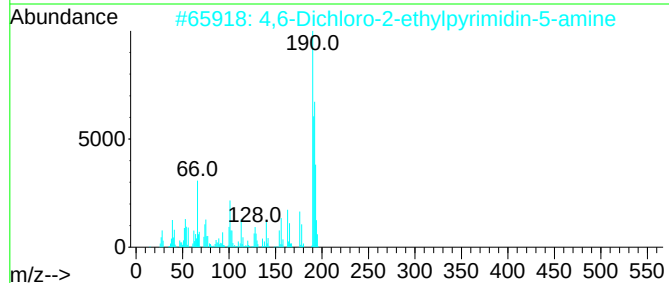
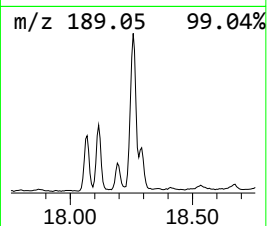
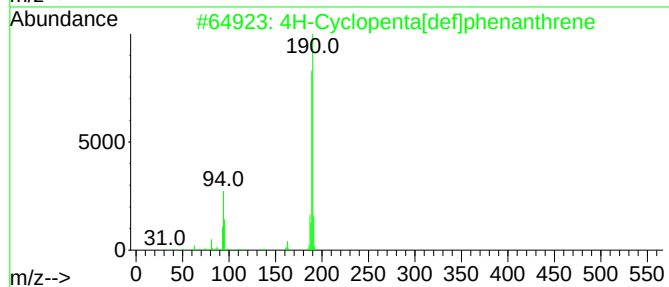
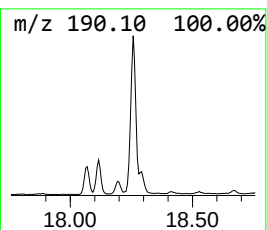
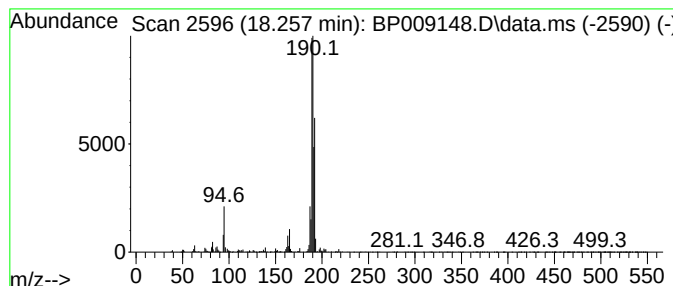
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 4H-Cyclopenta[def]phenanthrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.257	7.02 ng	1010220	Phenanthrene-d10	17.192

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	92
2			4,6-Dichloro-2-ethylpyrimidin-5-...	191	C6H7Cl2N3	006237-96-3	43
3			Benzaldehyde, 3,5-dichloro-2-hyd...	190	C7H4Cl2O2	000090-60-8	40
4			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	40
5			2-Amino-4,6-dichloropyrimidine-5...	191	C5H3Cl2N3O	005604-46-6	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021422\
Data File : BP009148.D
Acq On : 14 Feb 2022 22:47
Operator : CG/JU
Sample : N1526-03 2X
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ALS Vial : 19 Sample Multiplier: 1

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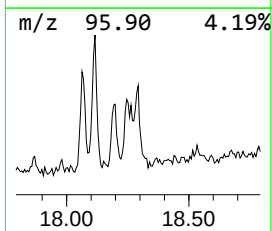
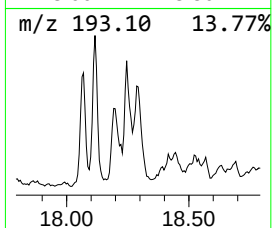
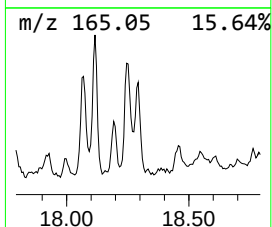
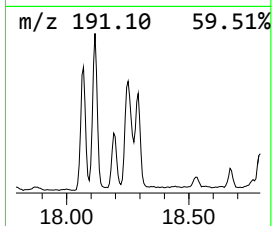
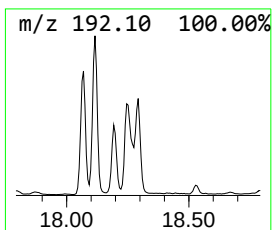
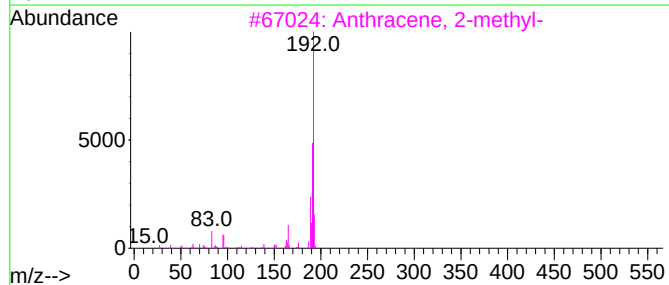
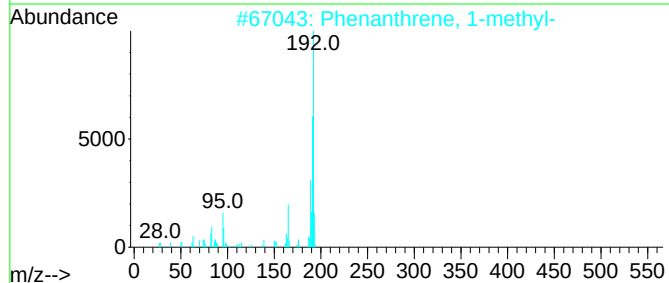
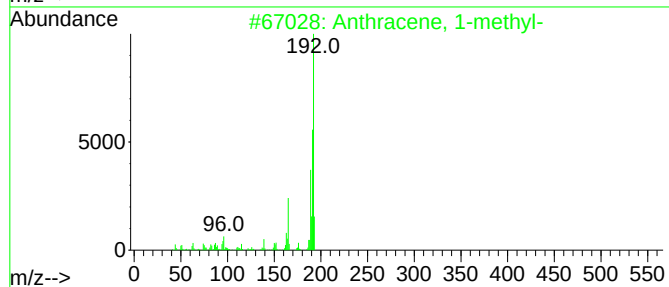
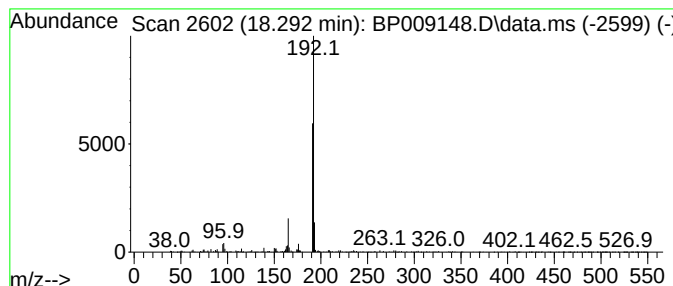
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Anthracene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.292	2.89 ng	416183	Phenanthrene-d10	17.192

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 1-methyl-	192	C15H12	000610-48-0	91
2		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	91
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
5		Anthracene, 9-methyl-	192	C15H12	000779-02-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021422\
Data File : BP009148.D
Acq On : 14 Feb 2022 22:47
Operator : CG/JU
Sample : N1526-03 2X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
HE-2-021122

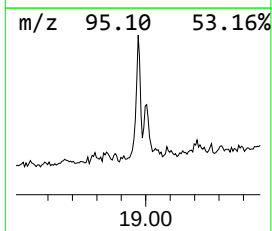
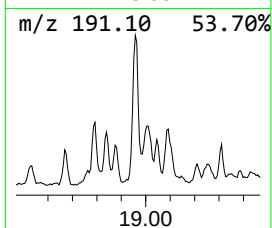
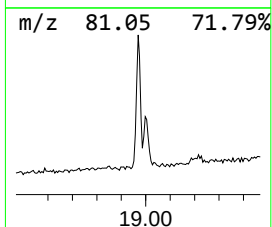
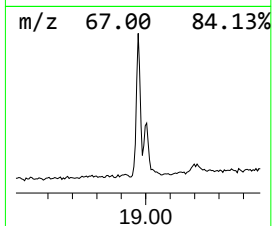
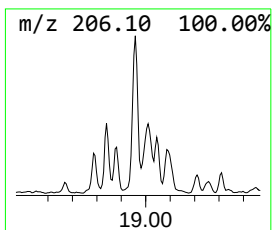
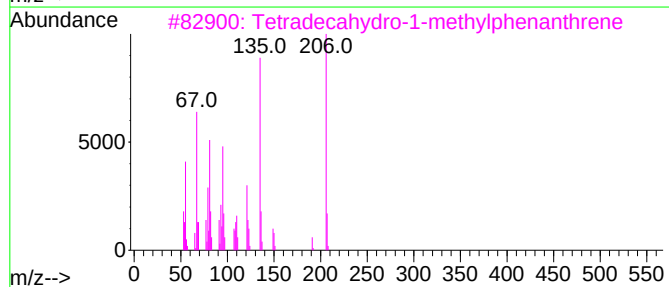
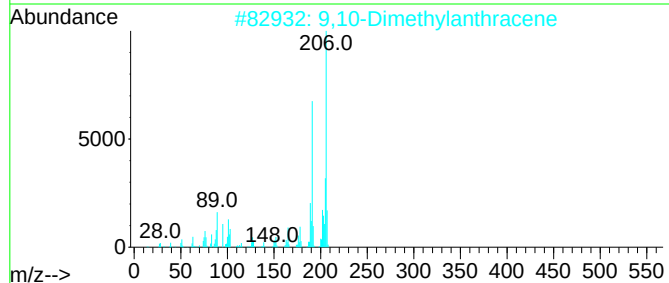
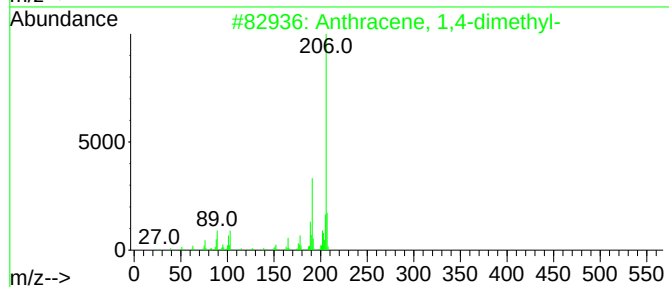
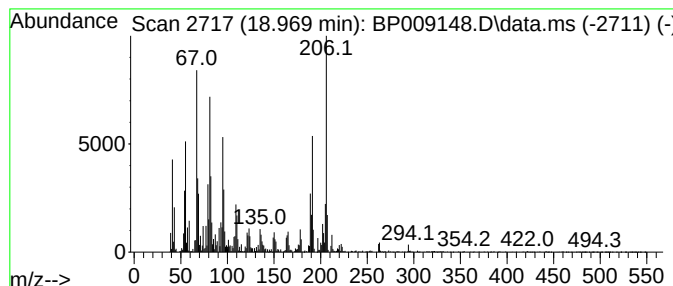
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Anthracene, 1,4-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.969	7.89 ng	1135580	Phenanthrene-d10	17.192

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	66
2		9,10-Dimethylantracene	206	C16H14	000781-43-1	64
3		Tetradecahydro-1-methylphenanthrene	206	C15H26	1000080-19-6	64
4		3,4-Dimethylphenanthrene	206	C16H14	1000452-24-5	64
5		Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021422\
Data File : BP009148.D
Acq On : 14 Feb 2022 22:47
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Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
HE-2-021122

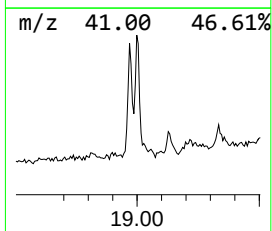
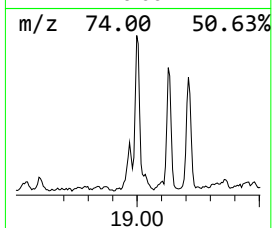
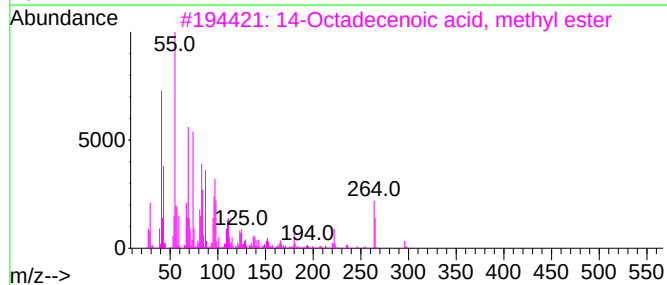
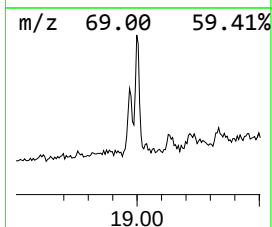
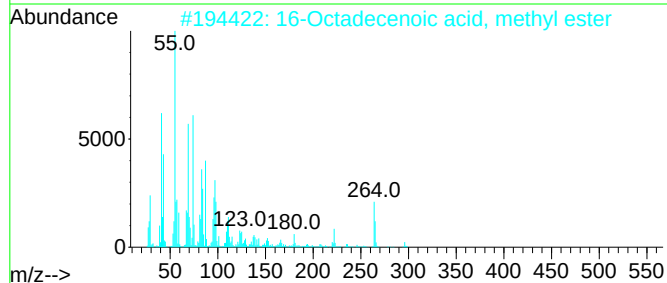
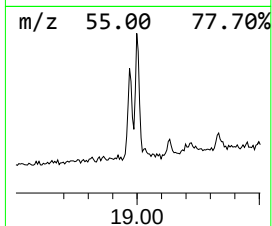
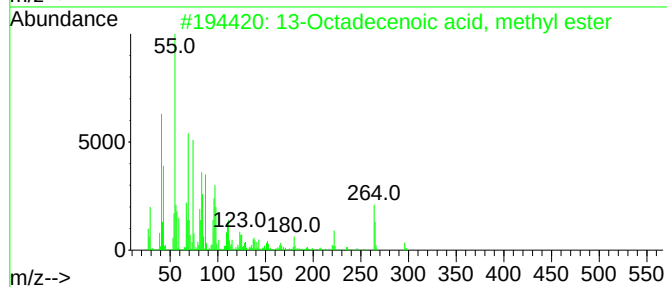
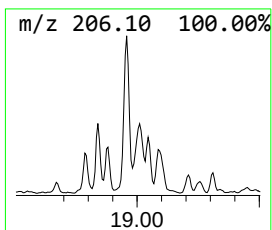
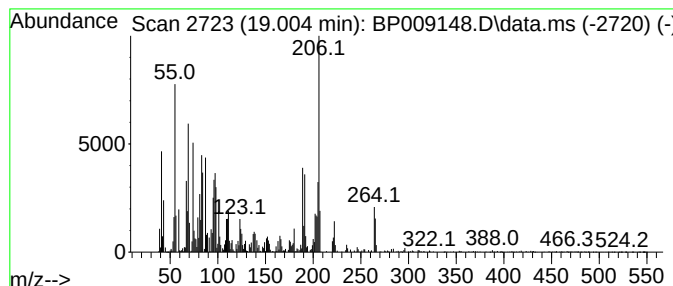
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 13-Octadecenoic acid, methyl... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.004	4.82 ng	693104	Phenanthrene-d10	17.192

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	94
2			16-Octadecenoic acid, methyl ester	296	C19H36O2	056554-49-5	94
3			14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	94
4			15-Octadecenoic acid, methyl ester	296	C19H36O2	004764-72-1	94
5			11-Octadecenoic acid, methyl est...	296	C19H36O2	001937-63-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021422\
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Misc :
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Instrument :
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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
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Dibenzothiophene	17.016	2.3	ng	327167	4	17.192	2877660	20.0
Phenanthrene, 2...	18.069	4.2	ng	608263	4	17.192	2877660	20.0
Anthracene, 2-m...	18.116	5.5	ng	786189	4	17.192	2877660	20.0
1H-Cyclopropa[1...	18.192	2.0	ng	292117	4	17.192	2877660	20.0
4H-Cyclopenta[d...	18.257	7.0	ng	1010220	4	17.192	2877660	20.0
Anthracene, 1-m...	18.292	2.9	ng	416183	4	17.192	2877660	20.0
Anthracene, 1,4...	18.969	7.9	ng	1135580	4	17.192	2877660	20.0
13-Octadecenoic...	19.004	4.8	ng	693104	4	17.192	2877660	20.0