

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
 Data File : BP007526.D
 Acq On : 20 Oct 2021 13:35
 Operator : CG/JU
 Sample : M4207-07
 Misc :
 ALS Vial : 81 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AY7

Integration Parameters: LSCINT.P

Integrator: RTE

Soothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % 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\SFAM-EPA-BP101421.M
 Title : SVOA CALIBRATION

Signal : TIC: BP007526.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.381	12	16	24	rVB	35428	49717	0.51%	0.044%
2	3.905	101	105	112	rVB	25650	34548	0.35%	0.031%
3	4.040	123	128	135	rVB2	55840	85355	0.87%	0.076%
4	4.134	140	144	149	rVB	23598	33930	0.35%	0.030%
5	4.228	156	160	167	rVB2	23203	35377	0.36%	0.032%
6	5.052	294	300	305	rBV	20107	32817	0.34%	0.029%
7	6.334	512	518	524	rBV2	21980	33210	0.34%	0.030%
8	7.110	642	650	660	rBV	457048	732686	7.50%	0.656%
9	7.281	672	679	687	rBV	380271	592867	6.07%	0.531%
10	7.481	707	713	724	rVB	470901	746266	7.64%	0.668%
11	7.952	787	793	801	rBV	917756	1460761	14.95%	1.307%
12	8.493	879	885	894	rVV2	25000	54177	0.55%	0.048%
13	8.657	906	913	922	rBV	577020	896776	9.18%	0.803%
14	8.775	928	933	940	rVB	28660	50499	0.52%	0.045%
15	9.110	982	990	1001	rVB	419641	692102	7.08%	0.619%
16	9.834	1105	1113	1123	rBV	294813	489638	5.01%	0.438%
17	10.375	1198	1205	1213	rBV	549302	889768	9.11%	0.796%
18	10.757	1259	1270	1276	rBV	1208206	2072241	21.21%	1.854%
19	10.810	1276	1279	1284	rVV	36325	57409	0.59%	0.051%
20	10.887	1284	1292	1305	rVV	474303	826510	8.46%	0.740%
21	12.245	1518	1523	1532	rVV	15622	37997	0.39%	0.034%
22	12.416	1546	1552	1561	rVB	40605	68424	0.70%	0.061%
23	12.634	1584	1589	1595	rVB	18959	30528	0.31%	0.027%
24	13.916	1802	1807	1811	rBV4	17378	30553	0.31%	0.027%
25	13.998	1811	1821	1833	rVV	969860	1345634	13.77%	1.204%
26	14.281	1860	1869	1884	rBV	1099213	1820397	18.63%	1.629%
27	14.586	1912	1921	1928	rBV	1810731	2741060	28.06%	2.453%
28	14.769	1946	1952	1962	rBV	298779	452740	4.63%	0.405%
29	14.986	1984	1989	1997	rBV2	39074	82931	0.85%	0.074%
30	15.216	2020	2028	2031	rBV5	11690	28813	0.29%	0.026%
31	15.257	2031	2035	2041	rVB5	18512	29824	0.31%	0.027%
32	15.386	2053	2057	2064	rVB7	18625	40346	0.41%	0.036%
33	15.581	2084	2090	2096	rBV	1447094	2067899	21.17%	1.851%
34	15.633	2096	2099	2104	rVV3	65165	94838	0.97%	0.085%
35	15.692	2104	2109	2114	rVB	176366	242284	2.48%	0.217%
36	15.933	2146	2150	2160	rVB5	23988	49922	0.51%	0.045%

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Integrator: RTE

Smoothing : OFF

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Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
 Title : SVOA CALIBRATION

37	16.081	2171	2175	2182	rVB2	30600	62175	0.64%	0.056%
38	16.310	2208	2214	2217	rBV4	44691	91985	0.94%	0.082%
39	16.351	2218	2221	2230	rVB2	47245	67278	0.69%	0.060%
40	16.916	2313	2317	2322	rVB2	74065	99178	1.02%	0.089%
41	16.992	2326	2330	2338	rVB5	20790	46995	0.48%	0.042%
42	17.169	2357	2360	2368	rVB4	33839	60293	0.62%	0.054%
43	17.339	2383	2389	2394	rBV	2225502	3330882	34.10%	2.981%
44	17.380	2394	2396	2401	rVV	379260	448357	4.59%	0.401%
45	17.439	2401	2406	2409	rVV2	1518845	2377367	24.34%	2.128%
46	17.475	2409	2412	2418	rVB	3321400	4407105	45.11%	3.944%
47	17.739	2453	2457	2461	rVB	134531	176700	1.81%	0.158%
48	17.863	2475	2478	2483	rVB4	36368	49898	0.51%	0.045%
49	18.022	2502	2505	2509	rVB5	37202	44284	0.45%	0.040%
50	18.210	2533	2537	2538	rBV	76347	88902	0.91%	0.080%
51	18.233	2538	2541	2542	rVV	129408	141619	1.45%	0.127%
52	18.257	2542	2545	2550	rVB	178665	233025	2.39%	0.209%
53	18.333	2553	2558	2563	rBV	340099	459821	4.71%	0.411%
54	18.398	2563	2569	2572	rBV2	139742	259801	2.66%	0.232%
55	18.669	2611	2615	2616	rBV	65399	79995	0.82%	0.072%
56	18.686	2616	2618	2622	rVV	103299	133063	1.36%	0.119%
57	18.722	2622	2624	2630	rVB2	47966	62874	0.64%	0.056%
58	18.927	2657	2659	2663	rVB	46376	48960	0.50%	0.044%
59	18.974	2665	2667	2670	rBV	38945	42515	0.44%	0.038%
60	19.010	2671	2673	2678	rVB	51886	66612	0.68%	0.060%
61	19.092	2682	2687	2691	rBV	140671	232573	2.38%	0.208%
62	19.157	2692	2698	2701	rBV4	83012	150462	1.54%	0.135%
63	19.227	2706	2710	2718	rVV3	123478	217259	2.22%	0.194%
64	19.345	2725	2730	2735	rBV	2015325	2682683	27.46%	2.401%
65	19.463	2748	2750	2753	rVB2	48262	40963	0.42%	0.037%
66	19.580	2768	2770	2773	rBV	47725	59923	0.61%	0.054%
67	19.669	2780	2785	2788	rBV2	2214435	3303470	33.81%	2.956%
68	19.698	2788	2790	2798	rVV	1891040	2288176	23.42%	2.048%
69	19.774	2799	2803	2810	rVB2	106611	200922	2.06%	0.180%
70	19.874	2816	2820	2826	rBV	266941	353259	3.62%	0.316%
71	20.021	2839	2845	2851	rBV3	263412	596634	6.11%	0.534%
72	20.180	2862	2872	2876	rBV	702491	1378487	14.11%	1.234%
73	20.280	2884	2889	2893	rBV	777946	952204	9.75%	0.852%
74	20.333	2893	2898	2903	rVV2	313949	512572	5.25%	0.459%
75	20.474	2918	2922	2925	rBV	343169	466521	4.78%	0.417%
76	20.874	2987	2990	2993	rBV2	100301	155676	1.59%	0.139%

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 Sampling : 1 Min Area: 0.1 % of largest Peak
 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

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77	20.910	2993	2996	3003	rVB	365273	563331	5.77%	0.504%
78	21.080	3018	3025	3028	rBV3	606345	1026206	10.50%	0.918%
79	21.116	3028	3031	3033	rVV	202384	242332	2.48%	0.217%
80	21.151	3033	3037	3042	rVB2	337049	504811	5.17%	0.452%
81	21.427	3076	3084	3087	rBV2	3298191	6998767	71.64%	6.263%
82	21.468	3087	3091	3100	rVB	6832775	9399196	96.21%	8.411%
83	21.604	3110	3114	3119	rVB3	235566	372237	3.81%	0.333%
84	21.698	3126	3130	3135	rVB	715773	1008120	10.32%	0.902%
85	21.751	3136	3139	3145	rVB3	272769	378281	3.87%	0.339%
86	22.015	3181	3184	3188	rVB	466376	615784	6.30%	0.551%
87	22.068	3189	3193	3199	rVB	562565	837014	8.57%	0.749%
88	22.227	3216	3220	3224	rBV	322474	472670	4.84%	0.423%
89	22.974	3344	3347	3353	rVB2	372646	599623	6.14%	0.537%
90	23.133	3367	3374	3388	rBV	2982669	9769279	100.00%	8.743%
91	23.321	3402	3406	3413	rVB	391268	664218	6.80%	0.594%
92	23.668	3458	3465	3470	rBV	2310454	4825933	49.40%	4.319%
93	23.721	3470	3474	3478	rVV2	1071287	1820424	18.63%	1.629%
94	23.774	3478	3483	3488	rVB	1415831	2552381	26.13%	2.284%
95	23.880	3494	3501	3506	rBV2	1667134	3610892	36.96%	3.231%
96	23.939	3506	3511	3520	rVB2	863058	2102718	21.52%	1.882%
97	24.574	3613	3619	3625	rBV	548235	1141675	11.69%	1.022%
98	26.451	3928	3938	3955	rVB2	2264852	8302262	84.98%	7.430%
99	26.856	4001	4007	4024	rVB	644090	2108147	21.58%	1.887%
100	27.239	4062	4072	4087	rVB	1886110	6325333	64.75%	5.661%

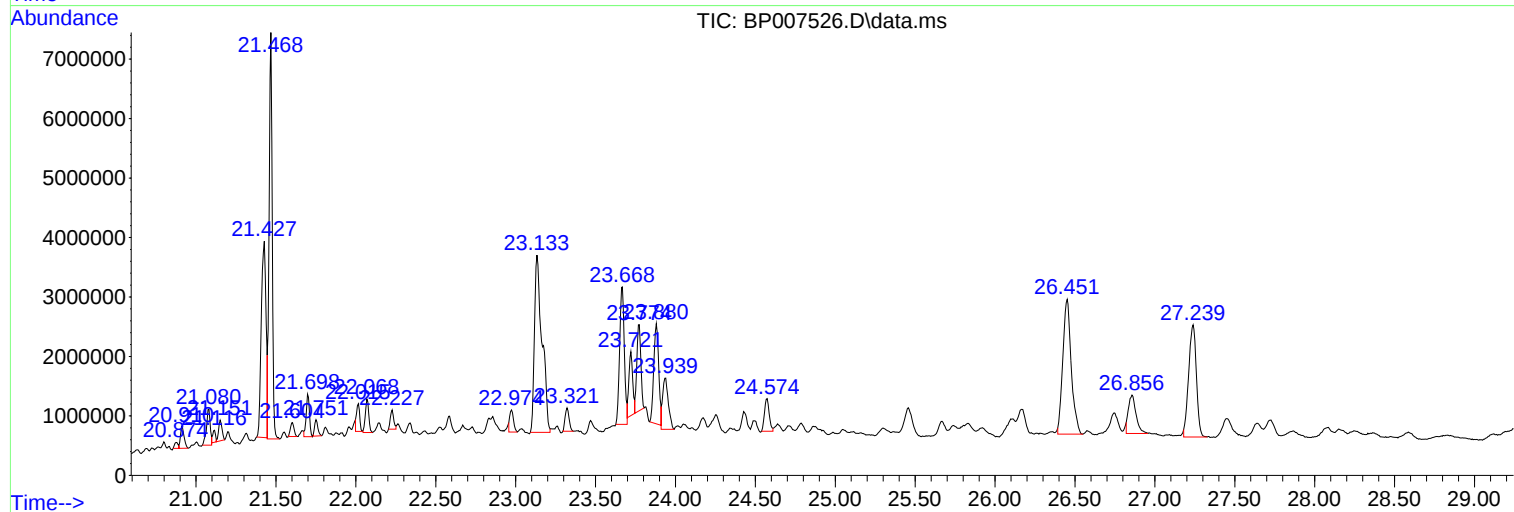
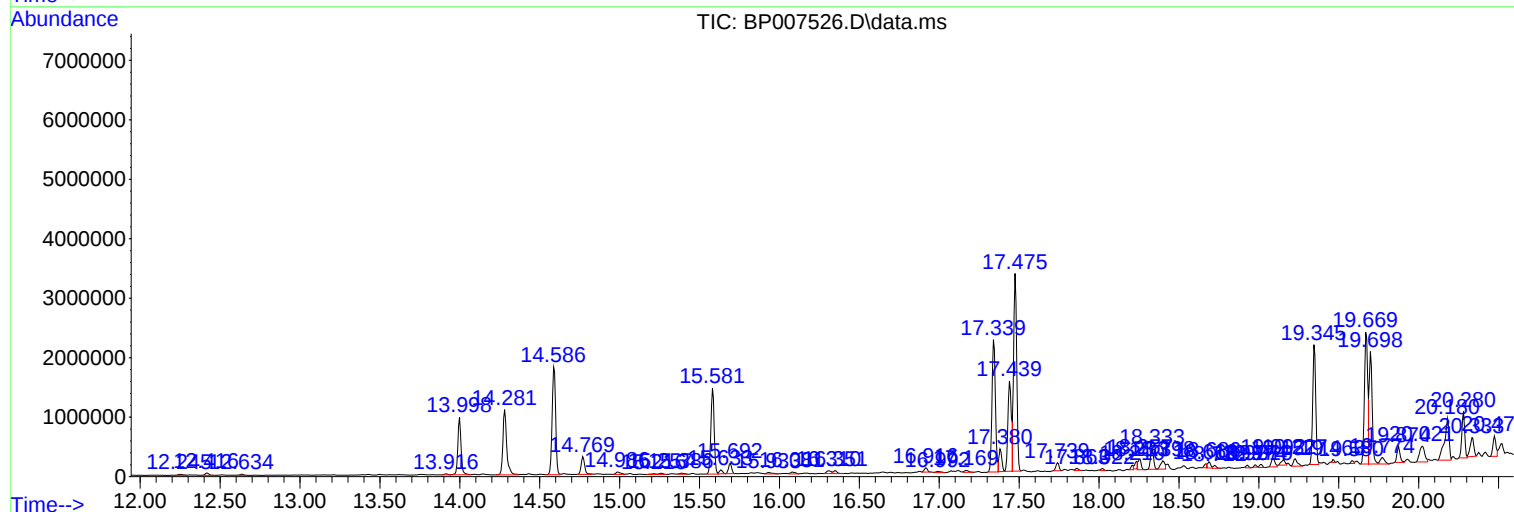
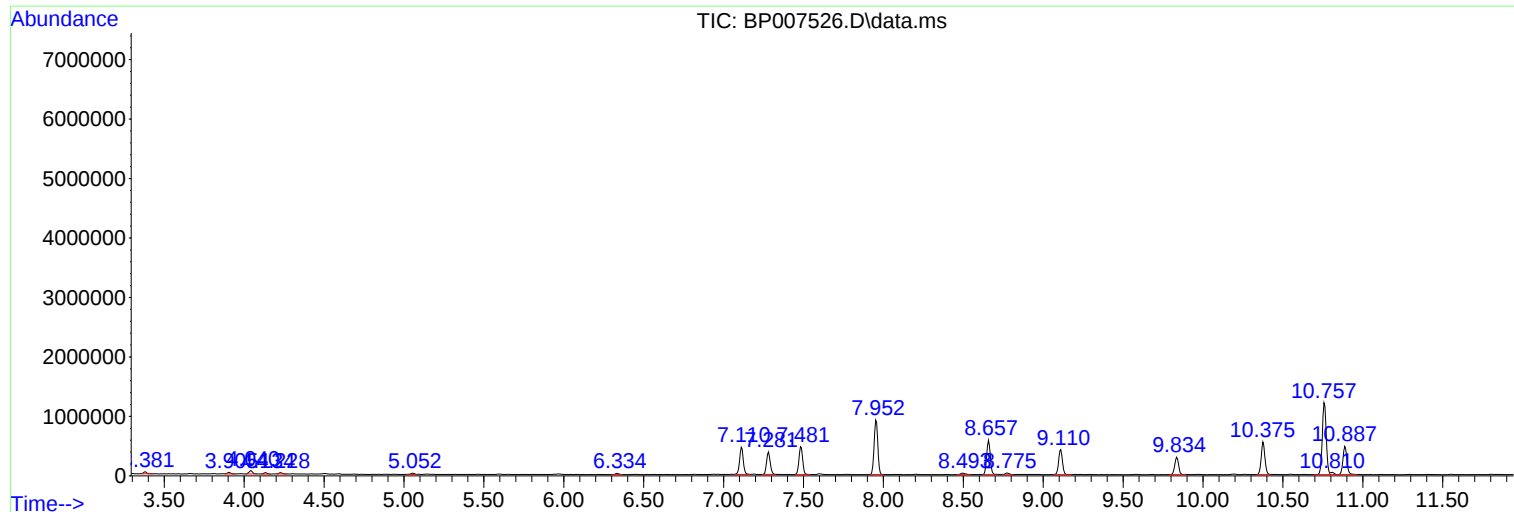
Sum of corrected areas: 111743946

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Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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



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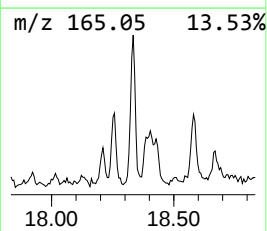
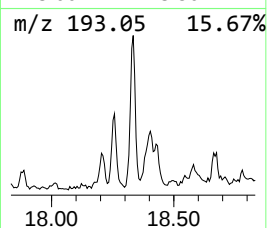
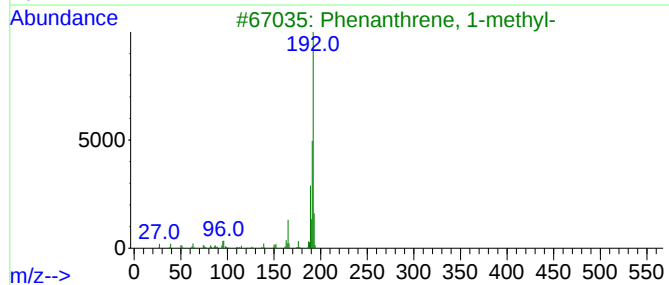
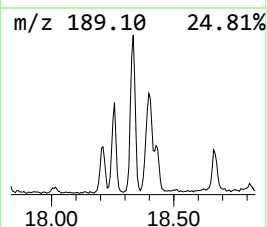
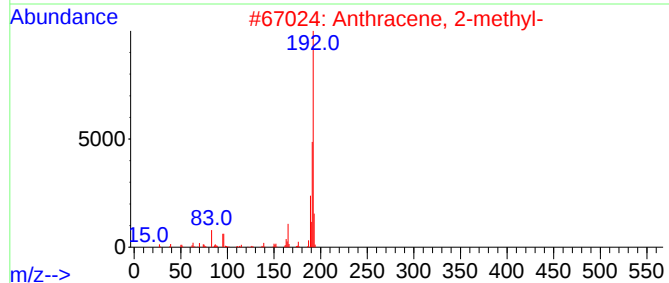
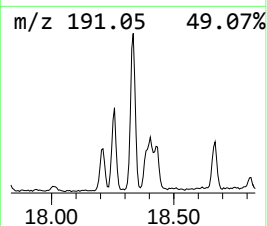
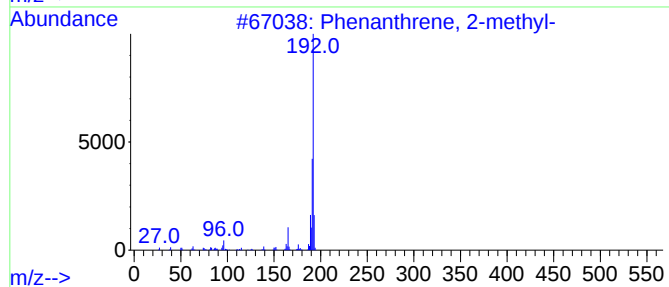
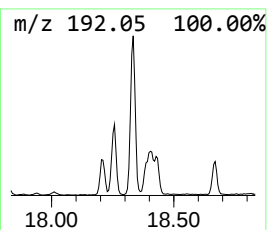
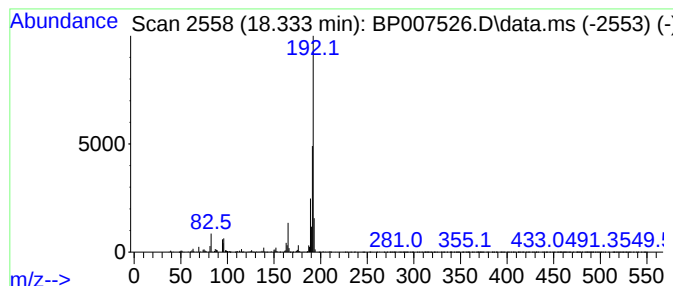
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Phenanthrene, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.333	2.76 ng/ul	459821	Phenanthrene-d10	17.339

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



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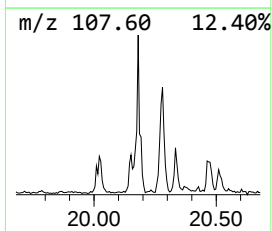
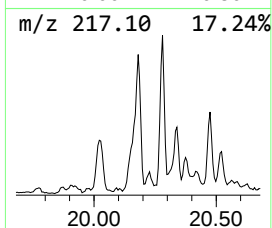
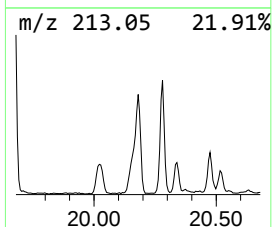
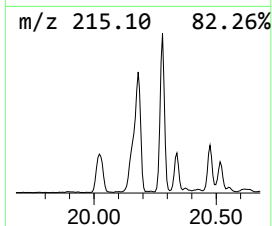
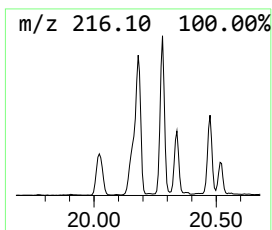
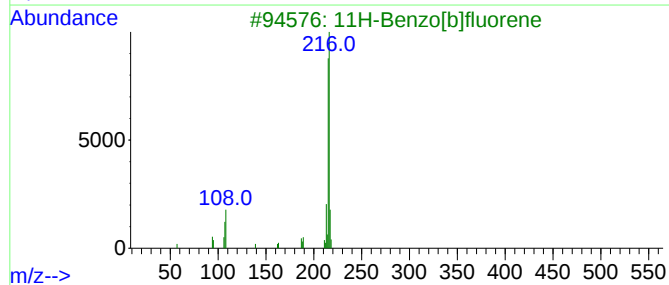
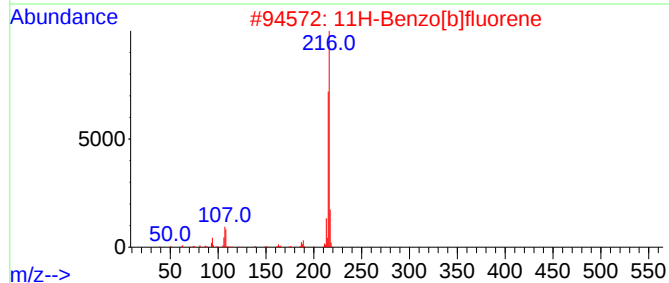
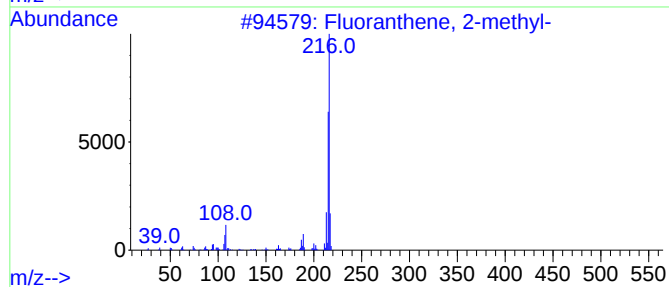
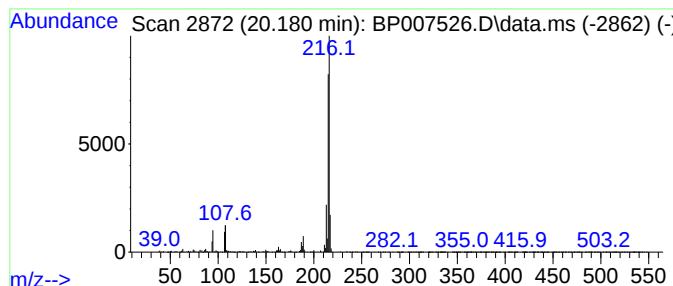
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Fluoranthene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.180	3.94 ng/ul	1378490	Chrysene-d12	21.427

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



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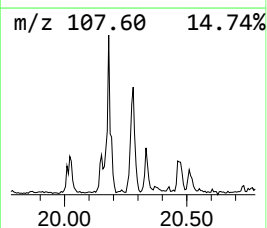
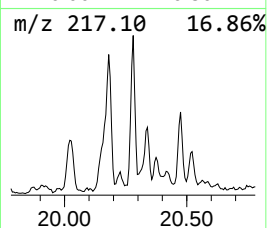
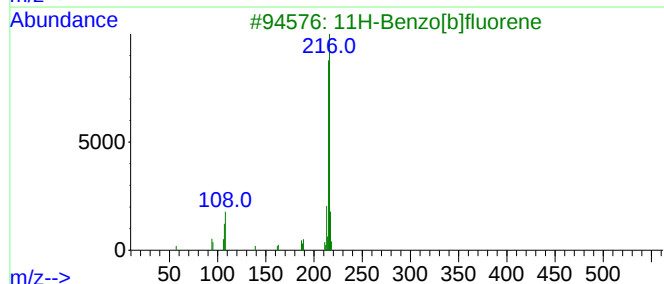
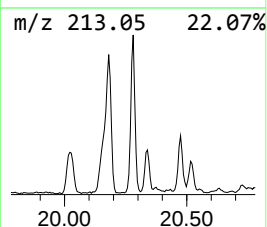
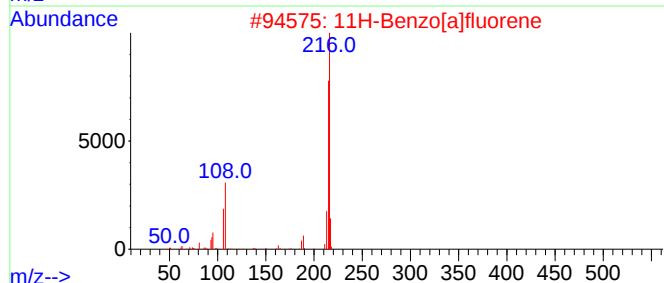
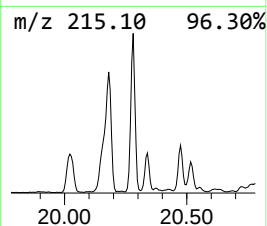
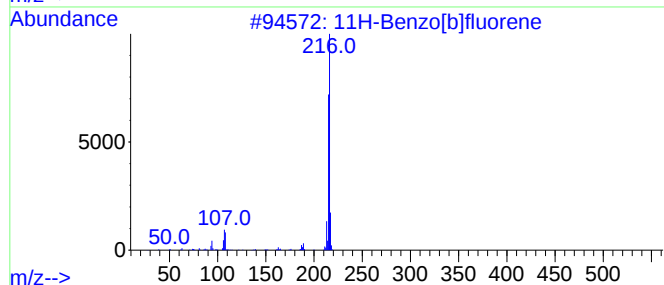
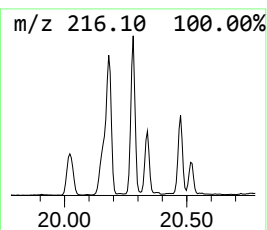
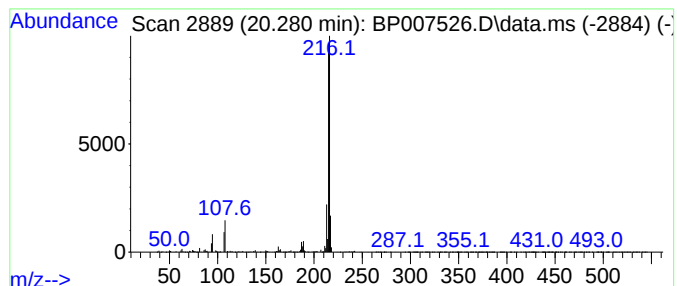
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 11H-Benzo[b]fluorene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.280	2.72 ng/ul	952204	Chrysene-d12	21.427

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	97
2			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	91
3			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
4			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
5			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007526.D
Acq On : 20 Oct 2021 13:35
Operator : CG/JU
Sample : M4207-07
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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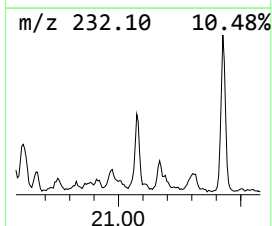
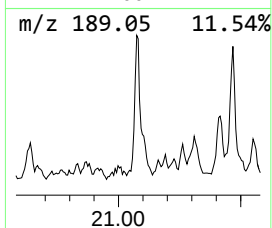
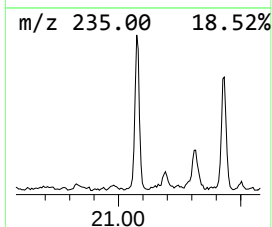
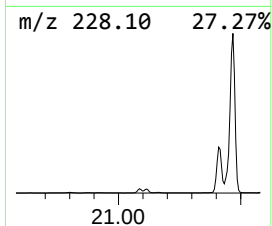
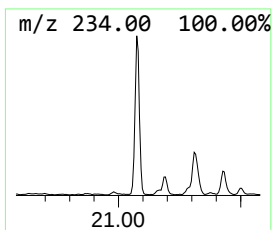
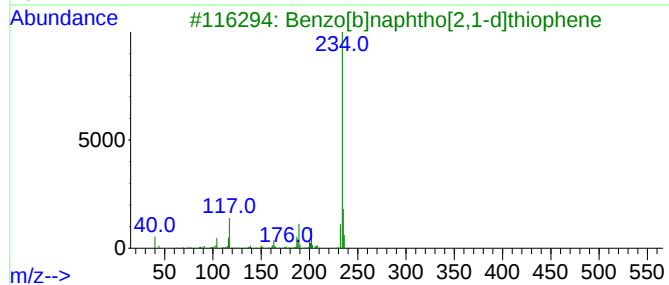
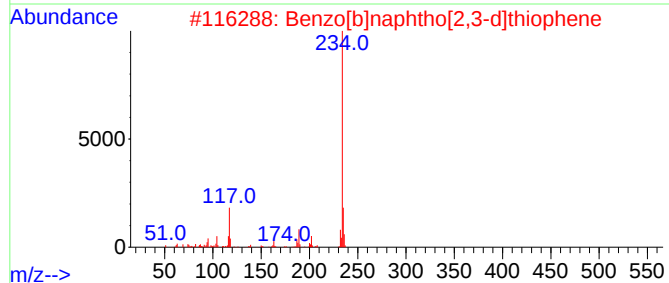
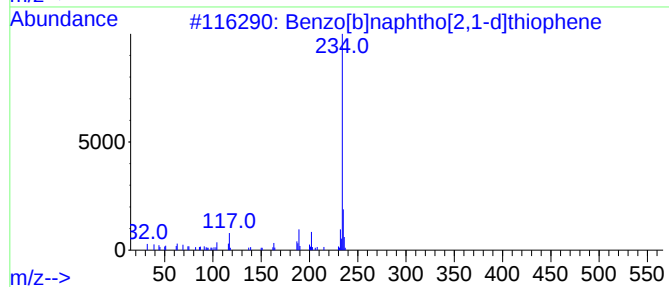
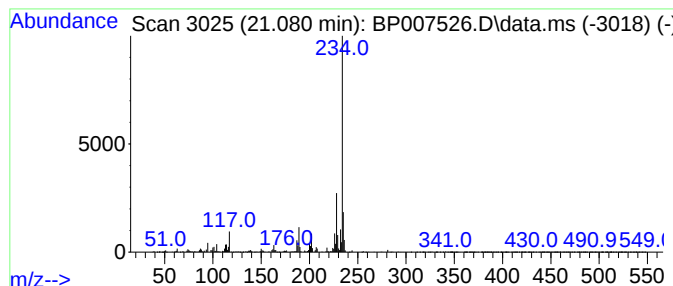
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzo[b]naphtho[2,1-d]thiop... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.080	2.93 ng/ul	1026210	Chrysene-d12	21.427

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	98
2			Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	95
3			Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	95
4			Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	94
5			Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007526.D
Acq On : 20 Oct 2021 13:35
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Sample : M4207-07
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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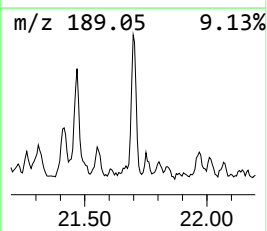
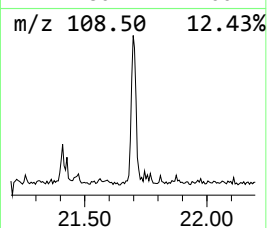
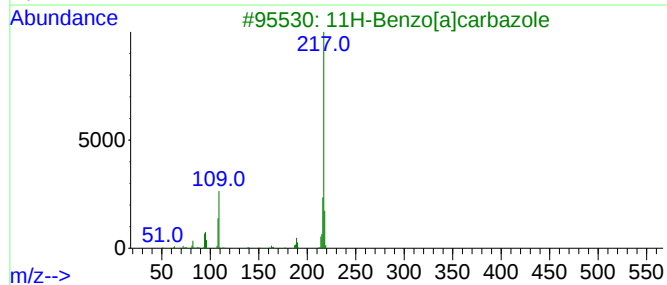
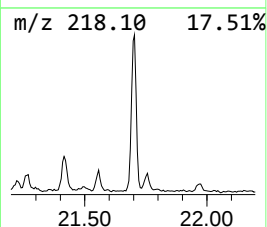
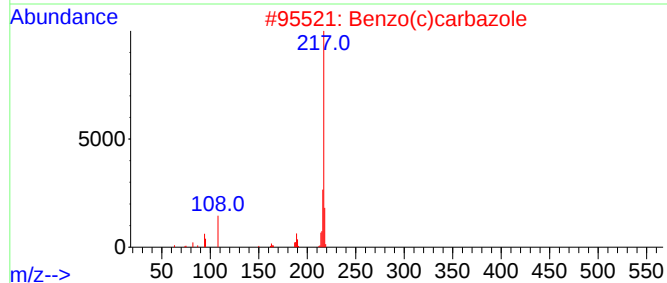
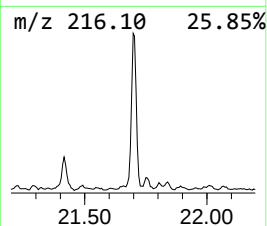
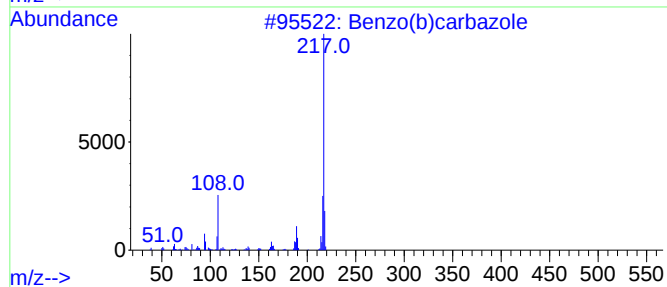
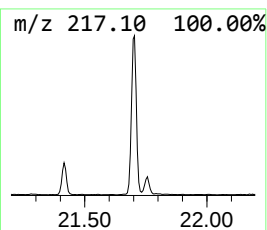
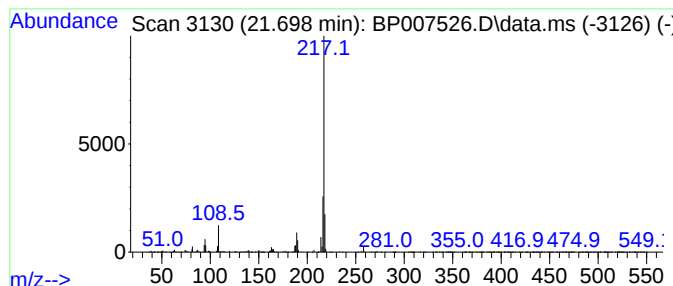
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzo(b)carbazole Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.698	2.88 ng/ul	1008120	Chrysene-d12	21.427

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo(b)carbazole	217	C16H11N	000243-28-7	96
2			Benzo(c)carbazole	217	C16H11N	034777-33-8	95
3			11H-Benzo[a]carbazole	217	C16H11N	000239-01-0	91
4			Benzo(b)carbazole	217	C16H11N	000243-28-7	91
5			11H-Benzo[a]carbazole	217	C16H11N	000239-01-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007526.D
Acq On : 20 Oct 2021 13:35
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Sample : M4207-07
Misc :
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Instrument :
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ClientSampleId :
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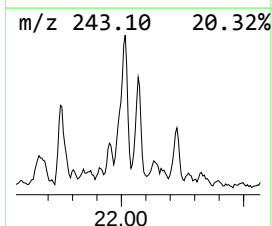
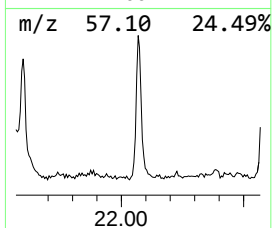
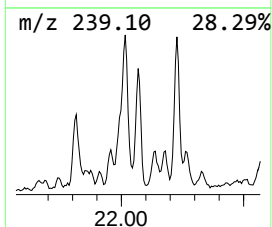
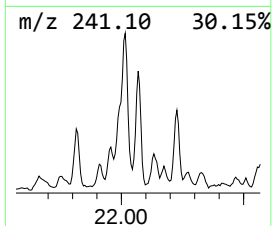
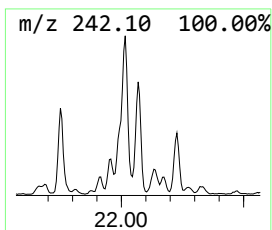
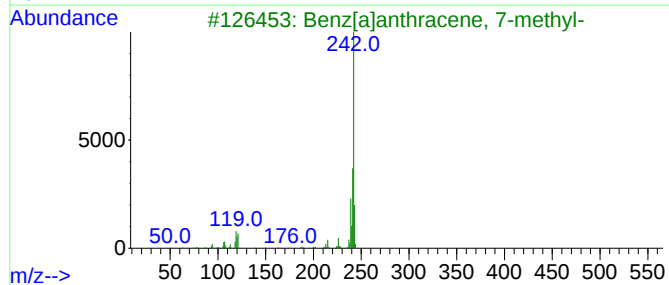
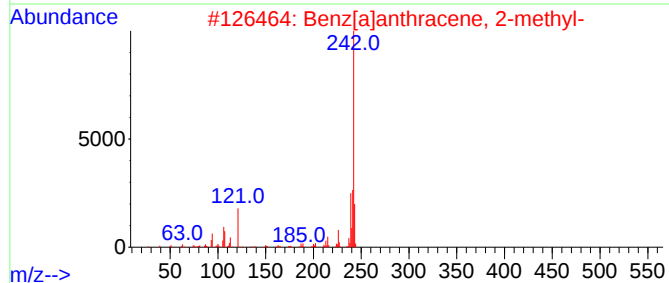
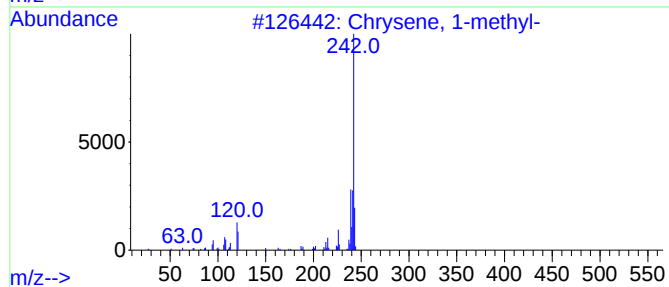
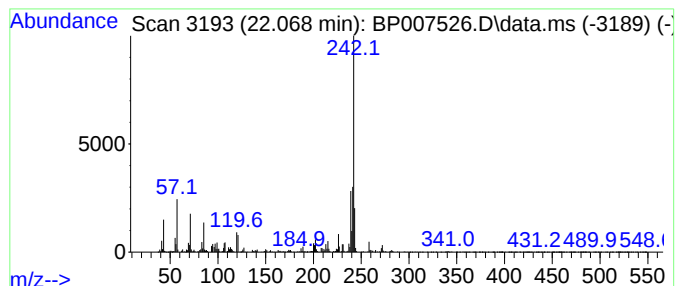
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Chrysene, 1-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.068	2.39 ng/ul	837014	Chrysene-d12	21.427

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Chrysene, 1-methyl-	242	C19H14	003351-28-8	98
2			Benz[a]anthracene, 2-methyl-	242	C19H14	002498-76-2	95
3			Benz[a]anthracene, 7-methyl-	242	C19H14	002541-69-7	94
4			1H-Benz[f]indene, 2-phenyl-	242	C19H14	1000305-22-4	93
5			Triphenylene, 2-methyl-	242	C19H14	001705-84-6	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007526.D
Acq On : 20 Oct 2021 13:35
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Sample : M4207-07
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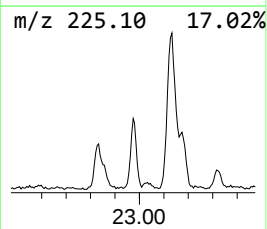
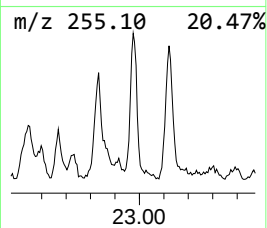
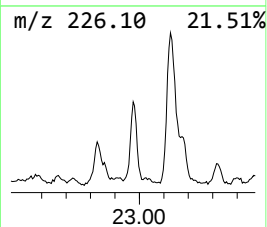
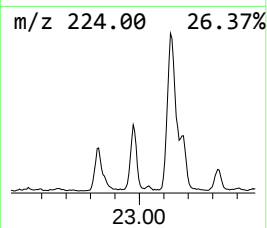
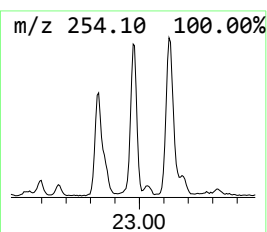
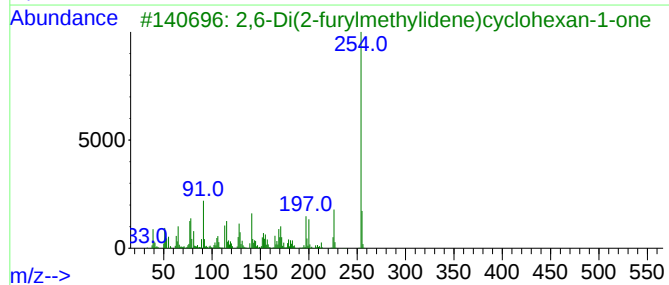
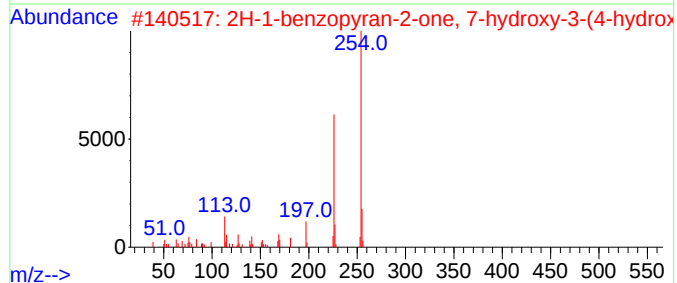
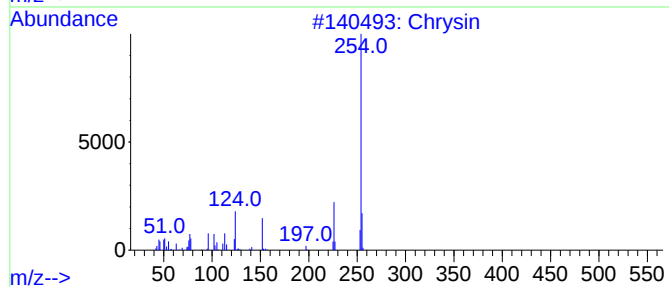
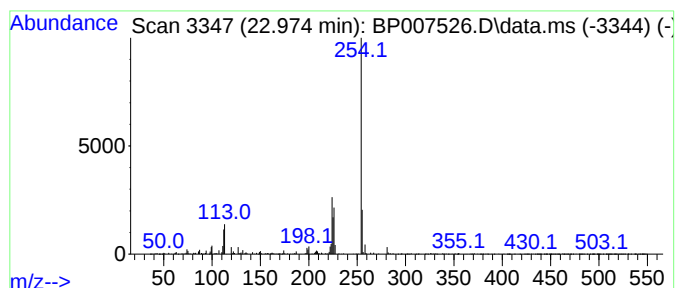
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Chrysin Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.974	3.32 ng/ul	599623	Perylene-d12	23.880

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Chrysin	254	C15H10O4	000480-40-0	58
2			2H-1-benzopyran-2-one, 7-hydroxy...	254	C15H10O4	1000401-42-6	53
3			2,6-Di(2-furylmethylidene)cyclohexan-1-one	254	C16H14O3	000893-00-5	53
4			Anthracene, 9-phenyl-	254	C20H14	000602-55-1	53
5			Benzo(a)pyrene, 7,8-dihydro-	254	C20H14	017573-23-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
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Misc :
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Instrument :
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ClientSampleId :
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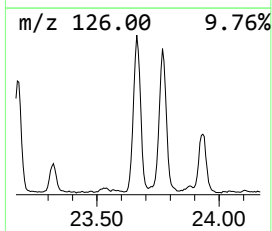
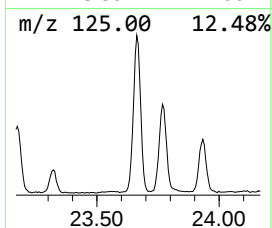
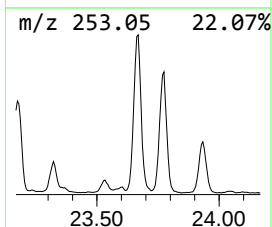
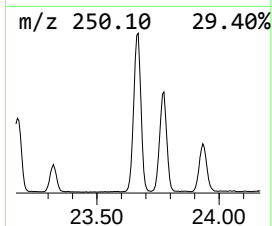
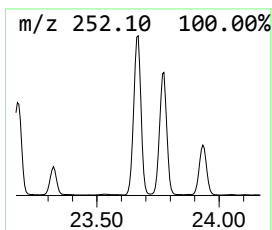
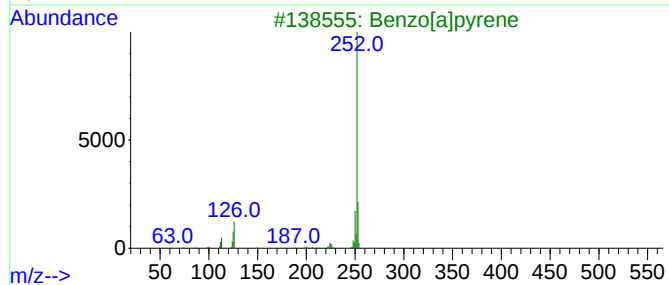
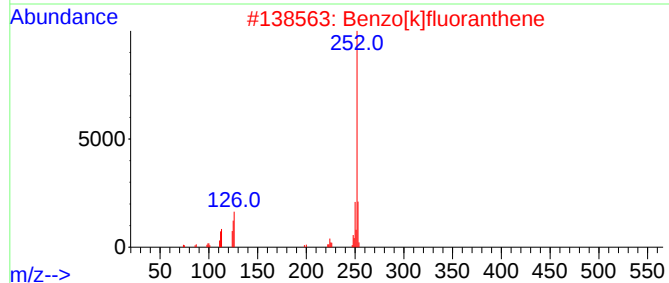
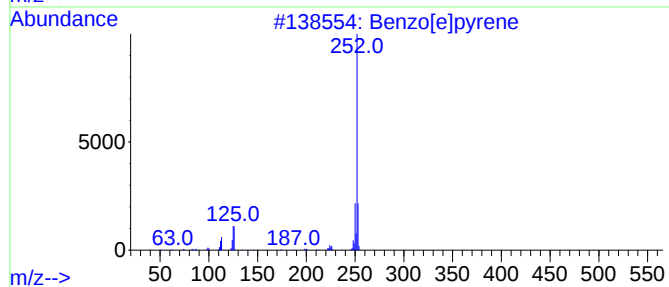
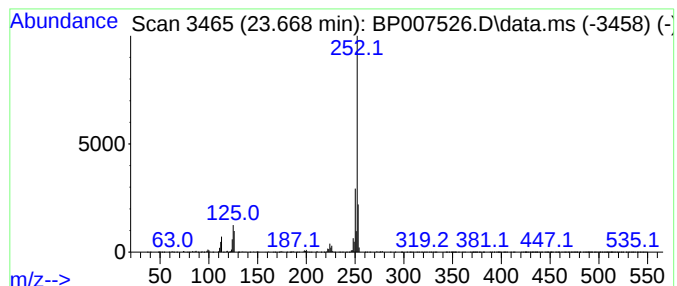
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzo[e]pyrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.668	26.73 ng/ul	4825930	Perylene-d12	23.880

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	99
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
3			Benzo[a]pyrene	252	C20H12	000050-32-8	94
4			Benzo[b]fluoranthene	252	C20H12	000205-99-2	94
5			Perylene	252	C20H12	000198-55-0	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
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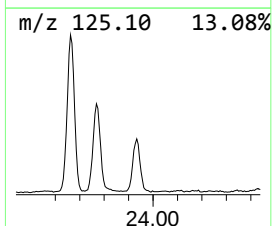
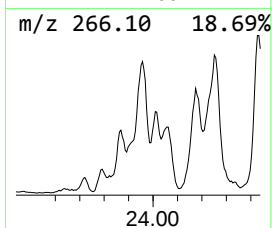
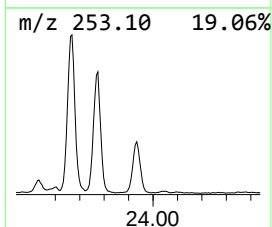
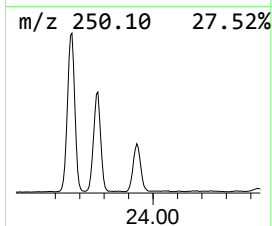
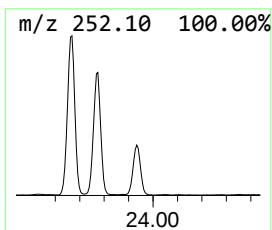
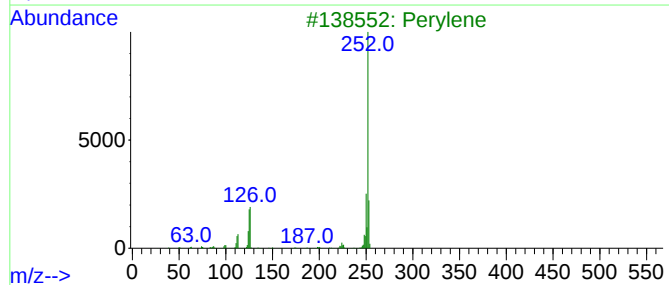
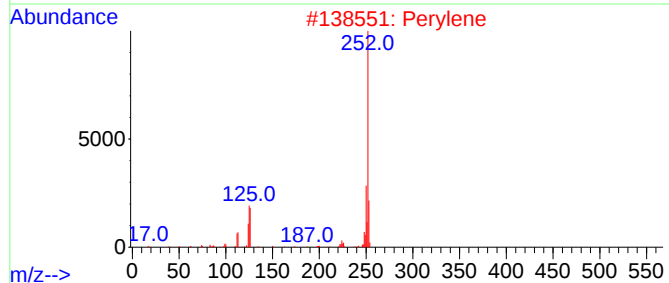
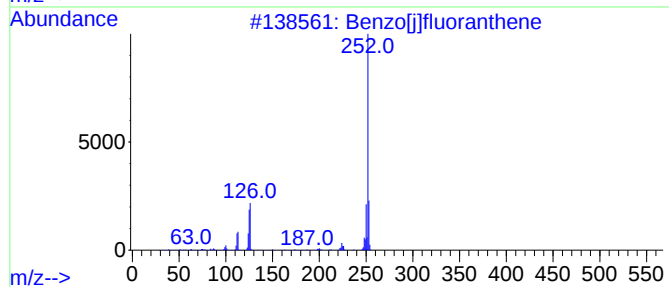
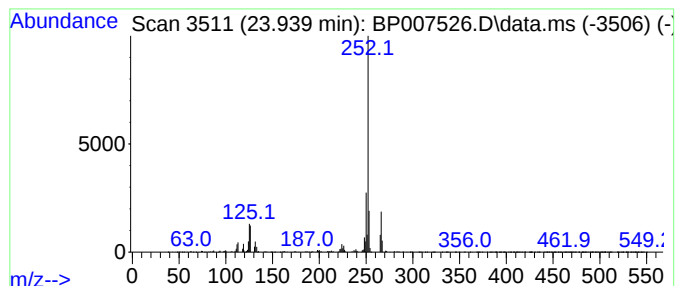
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Benzo[j]fluoranthene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.939	11.65 ng/ul	2102720	Perylene-d12	23.880

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[j]fluoranthene	252	C20H12	000205-82-3	96
2			Perylene	252	C20H12	000198-55-0	96
3			Perylene	252	C20H12	000198-55-0	96
4			Benzo[e]pyrene	252	C20H12	000192-97-2	96
5			Benzo[k]fluoranthene	252	C20H12	000207-08-9	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007526.D
Acq On : 20 Oct 2021 13:35
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Sample : M4207-07
Misc :
ALS Vial : 81 Sample Multiplier: 1

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ClientSampleId :
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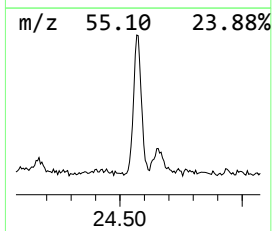
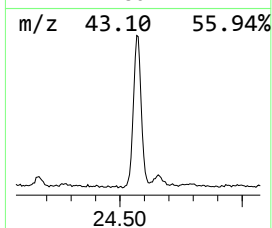
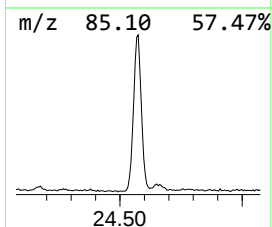
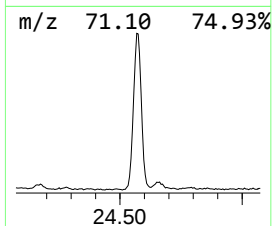
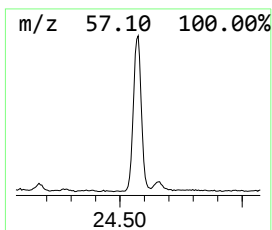
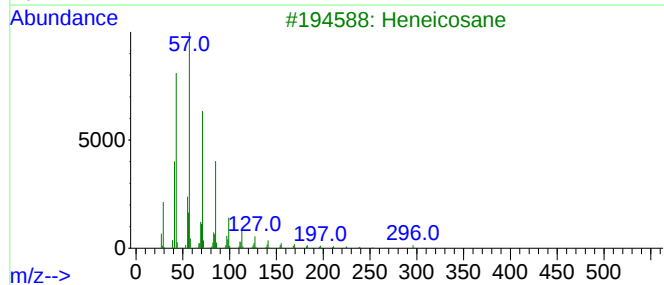
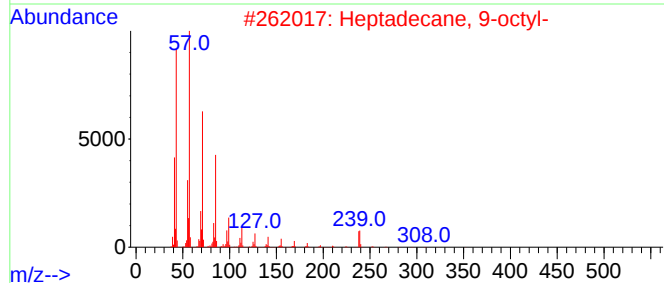
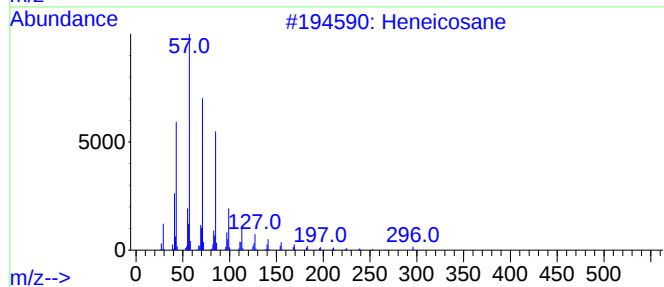
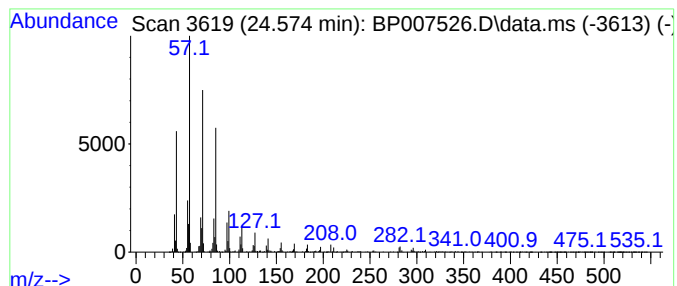
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.574	6.32 ng/ul	1141680	Perylene-d12	23.880

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	99
2			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	95
3			Heneicosane	296	C21H44	000629-94-7	93
4			Heneicosane, 11-decyl-	437	C31H64	055320-06-4	93
5			Eicosane	282	C20H42	000112-95-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
Data File : BP007526.D
Acq On : 20 Oct 2021 13:35
Operator : CG/JU
Sample : M4207-07
Misc :
ALS Vial : 81 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AY7

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
Quant Title : SVOA 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
Phenanthrene, 2...	18.333	2.8	ng/u1	459821	4	17.339	3330880	20.0
Fluoranthene, 2...	20.180	3.9	ng/u1	1378490	5	21.427	6998770	20.0
11H-Benzo[b]flu...	20.280	2.7	ng/u1	952204	5	21.427	6998770	20.0
Benzo[b]naphtho...	21.080	2.9	ng/u1	1026210	5	21.427	6998770	20.0
Benzo(b)carbazole	21.698	2.9	ng/u1	1008120	5	21.427	6998770	20.0
Chrysene, 1-met...	22.068	2.4	ng/u1	837014	5	21.427	6998770	20.0
Chrysin	22.974	3.3	ng/u1	599623	6	23.880	3610890	20.0
Benzo[e]pyrene	23.668	26.7	ng/u1	4825930	6	23.880	3610890	20.0
Benzo[j]fluoran...	23.939	11.7	ng/u1	2102720	6	23.880	3610890	20.0
(DEL) Alkane: S...	24.574	6.3	ng/u1	1141680	6	23.880	3610890	20.0