

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\
 Data File : BP004310.D
 Acq On : 15 Dec 2020 18:34
 Operator : CG/JU
 Sample : L5001-13DL 10X
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A54D1DL

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	7.004	645	649	658	rVB2	56245	111518	1.33%	0.145%
2	7.169	674	677	685	rVB	40385	68841	0.82%	0.089%
3	7.375	707	712	718	rVB	58515	95547	1.14%	0.124%
4	7.845	785	792	800	rVB	933795	1531121	18.25%	1.985%
5	8.545	906	911	919	rVB2	66314	106340	1.27%	0.138%
6	8.998	982	988	994	rVB	50006	87420	1.04%	0.113%
7	9.722	1104	1111	1117	rBV	40716	72851	0.87%	0.094%
8	10.263	1197	1203	1209	rVB	67110	113166	1.35%	0.147%
9	10.639	1259	1267	1273	rBV	1346177	2228188	26.55%	2.889%
10	10.692	1273	1276	1284	rVV	58460	93218	1.11%	0.121%
11	10.775	1284	1290	1301	rVB2	38472	74572	0.89%	0.097%
12	11.475	1403	1409	1415	rVB5	8349	17678	0.21%	0.023%
13	12.163	1517	1526	1531	rBV5	14485	36125	0.43%	0.047%
14	12.304	1544	1550	1557	rBV	45408	71555	0.85%	0.093%
15	12.528	1582	1588	1594	rVB	38207	67443	0.80%	0.087%
16	13.322	1718	1723	1729	rVB2	23405	37491	0.45%	0.049%
17	13.522	1751	1757	1766	rVB	13289	30177	0.36%	0.039%
18	13.651	1772	1779	1780	rBV3	23047	32095	0.38%	0.042%
19	13.810	1799	1806	1811	rBV	52510	97283	1.16%	0.126%
20	13.863	1811	1815	1816	rVV	34251	39524	0.47%	0.051%
21	13.892	1816	1820	1825	rVV	120535	178181	2.12%	0.231%
22	13.945	1825	1829	1833	rVV2	22163	32892	0.39%	0.043%
23	14.045	1842	1846	1850	rVV	27536	47033	0.56%	0.061%
24	14.180	1859	1869	1873	rVV	155890	261955	3.12%	0.340%
25	14.210	1873	1874	1882	rVV	77612	92564	1.10%	0.120%
26	14.322	1886	1893	1901	rVB5	8837	20885	0.25%	0.027%
27	14.492	1913	1922	1928	rBV	2040258	3082967	36.74%	3.997%
28	14.557	1928	1933	1940	rVV	294526	450234	5.37%	0.584%
29	14.622	1941	1944	1950	rVB2	21265	30655	0.37%	0.040%
30	14.692	1950	1956	1966	rBV3	49457	98955	1.18%	0.128%
31	14.804	1966	1975	1979	rBV4	25098	50508	0.60%	0.065%
32	14.892	1984	1990	1998	rVV	212079	313243	3.73%	0.406%
33	14.969	1998	2003	2008	rVV	34937	47253	0.56%	0.061%
34	15.110	2021	2027	2033	rBV2	36357	65619	0.78%	0.085%

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35	15.169	2033	2037	2041	rVV2	31632	44105	0.53%	0.057%
36	15.286	2049	2057	2060	rBV2	32003	66436	0.79%	0.086%
37	15.316	2060	2062	2067	rVV4	27733	37702	0.45%	0.049%
38	15.486	2083	2091	2096	rVV	197026	313006	3.73%	0.406%
39	15.545	2096	2101	2110	rVV2	382624	577019	6.88%	0.748%
40	15.639	2110	2117	2119	rVV3	36556	82194	0.98%	0.107%
41	15.669	2119	2122	2125	rVV3	59091	86814	1.03%	0.113%
42	15.710	2125	2129	2133	rVV3	60783	99183	1.18%	0.129%
43	15.757	2133	2137	2140	rVV3	35752	56295	0.67%	0.073%
44	15.798	2140	2144	2147	rVV4	32380	58630	0.70%	0.076%
45	15.839	2147	2151	2158	rVV	122197	205353	2.45%	0.266%
46	15.939	2161	2168	2169	rVV6	8788	19115	0.23%	0.025%
47	15.992	2171	2177	2184	rVV	134038	288339	3.44%	0.374%
48	16.080	2184	2192	2198	rVB3	34731	76644	0.91%	0.099%
49	16.221	2212	2216	2223	rVB5	40211	66449	0.79%	0.086%
50	16.363	2237	2240	2244	rBV4	11654	16297	0.19%	0.021%
51	16.557	2261	2273	2278	rBV3	106597	284983	3.40%	0.370%
52	16.604	2278	2281	2287	rVV2	63666	99793	1.19%	0.129%
53	16.657	2287	2290	2293	rVV5	23838	39878	0.48%	0.052%
54	16.704	2295	2298	2303	rVB2	42115	63876	0.76%	0.083%
55	16.786	2309	2312	2316	rBV2	56070	71138	0.85%	0.092%
56	16.827	2316	2319	2322	rVB2	68133	76225	0.91%	0.099%
57	16.904	2327	2332	2340	rVB	174715	275543	3.28%	0.357%
58	16.986	2340	2346	2356	rBV3	103905	315811	3.76%	0.409%
59	17.069	2356	2360	2370	rVB	205000	324264	3.86%	0.420%
60	17.257	2385	2392	2395	rBV	2498968	3787776	45.14%	4.911%
61	17.298	2395	2399	2404	rVV	4007563	5662184	67.48%	7.342%
62	17.351	2404	2408	2410	rVV2	274475	377144	4.49%	0.489%
63	17.386	2410	2414	2419	rVV	1032894	1426866	17.00%	1.850%
64	17.445	2419	2424	2430	rVB3	111880	184397	2.20%	0.239%
65	17.533	2435	2439	2442	rBV3	30614	55082	0.66%	0.071%
66	17.657	2451	2460	2464	rBV2	370884	632435	7.54%	0.820%
67	17.774	2477	2480	2484	rVB2	72677	77348	0.92%	0.100%
68	17.833	2487	2490	2495	rVV2	71819	89609	1.07%	0.116%
69	18.045	2523	2526	2530	rBV	46064	59862	0.71%	0.078%
70	18.127	2535	2540	2543	rVV	551603	841381	10.03%	1.091%
71	18.168	2543	2547	2553	rVV	763230	1098372	13.09%	1.424%

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72	18.251	2556	2561	2565	rVV	307211	431444	5.14%	0.559%
73	18.315	2566	2572	2575	rVV3	797213	1399911	16.68%	1.815%
74	18.345	2575	2577	2583	rVB	361965	469549	5.60%	0.609%
75	18.457	2594	2596	2600	rVB2	58398	60044	0.72%	0.078%
76	18.610	2617	2622	2625	rBV	451136	601758	7.17%	0.780%
77	18.651	2625	2629	2633	rVB	272091	373905	4.46%	0.485%
78	18.727	2640	2642	2648	rBV2	53353	64219	0.77%	0.083%
79	18.845	2659	2662	2666	rBV	139580	171533	2.04%	0.222%
80	18.898	2668	2671	2674	rVV	123631	151651	1.81%	0.197%
81	18.933	2674	2677	2680	rVB2	121880	136457	1.63%	0.177%
82	19.015	2686	2691	2695	rBV	287135	498080	5.94%	0.646%
83	19.151	2710	2714	2722	rVB2	254420	425372	5.07%	0.552%
84	19.268	2729	2734	2740	rBV	6331980	8324440	99.21%	10.794%
85	19.362	2747	2750	2754	rVB	150846	186277	2.22%	0.242%
86	19.592	2784	2789	2791	rBV3	1377141	1986472	23.67%	2.576%
87	19.621	2791	2794	2802	rVV	4873440	6892662	82.14%	8.937%
88	19.692	2803	2806	2812	rVB2	286257	428799	5.11%	0.556%
89	19.798	2821	2824	2830	rVB	328777	402615	4.80%	0.522%
90	19.857	2831	2834	2839	rVB	296263	378612	4.51%	0.491%
91	19.933	2843	2847	2848	rBV2	330607	425562	5.07%	0.552%
92	20.104	2873	2876	2881	rVB	672266	885249	10.55%	1.148%
93	20.204	2890	2893	2896	rBV	400126	442984	5.28%	0.574%
94	20.839	2998	3001	3006	rVB	549623	700308	8.35%	0.908%
95	21.127	3047	3050	3056	rVB2	698342	897858	10.70%	1.164%
96	21.345	3081	3087	3091	rBV3	4248038	8390953	100.00%	10.880%
97	21.386	3091	3094	3103	rVB	2463474	3207439	38.22%	4.159%
98	23.003	3364	3369	3374	rBV	1753924	3984804	47.49%	5.167%
99	23.615	3469	3473	3480	rVB	1806707	3449435	41.11%	4.473%
100	23.721	3486	3491	3497	rBV	1962474	3632617	43.29%	4.710%

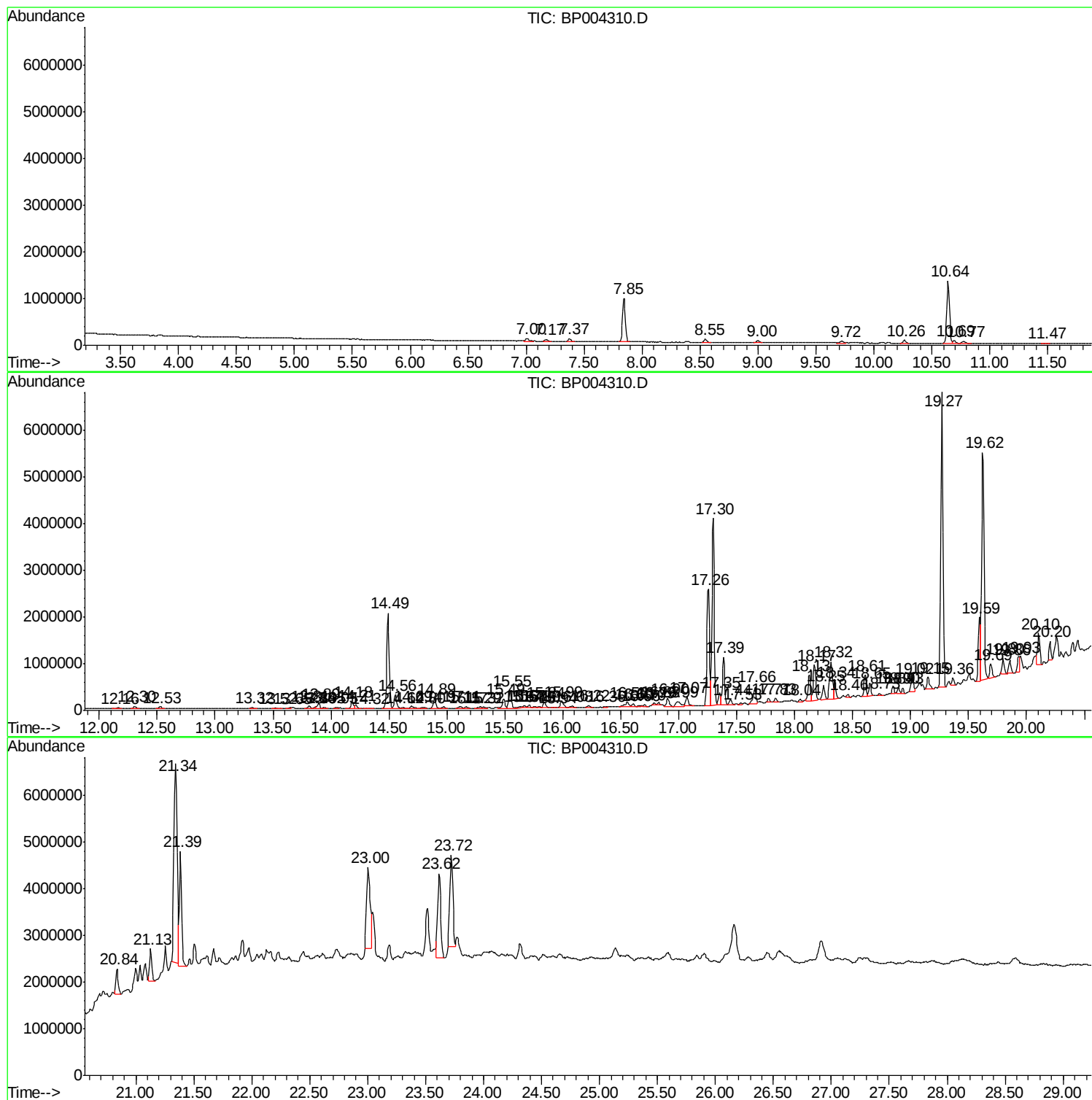
Sum of corrected areas: 77123654

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Instrument :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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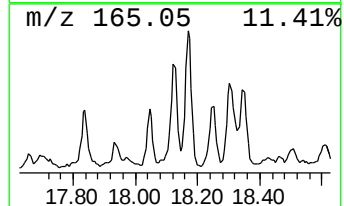
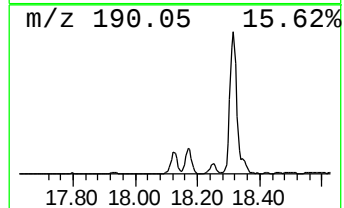
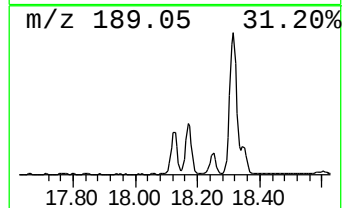
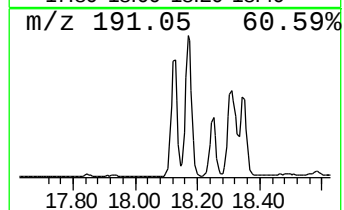
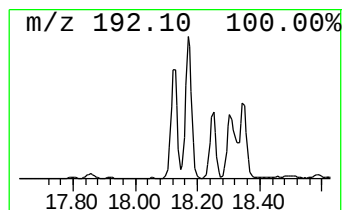
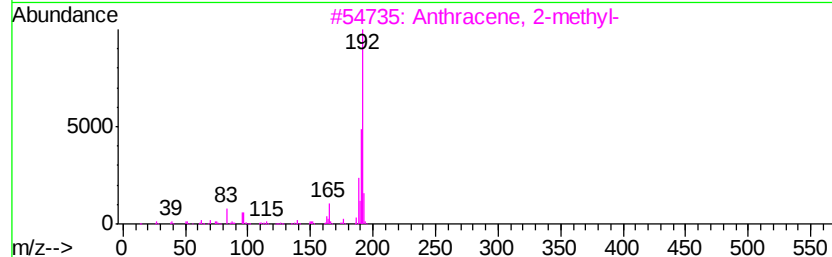
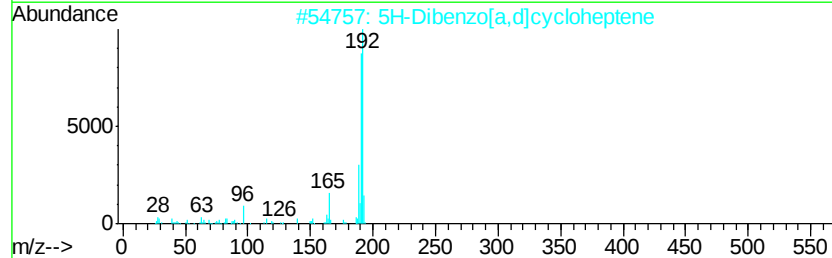
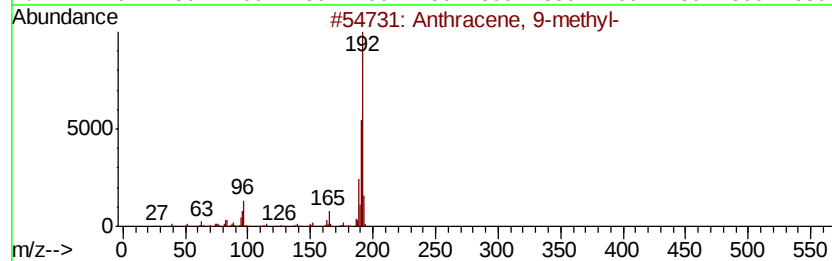
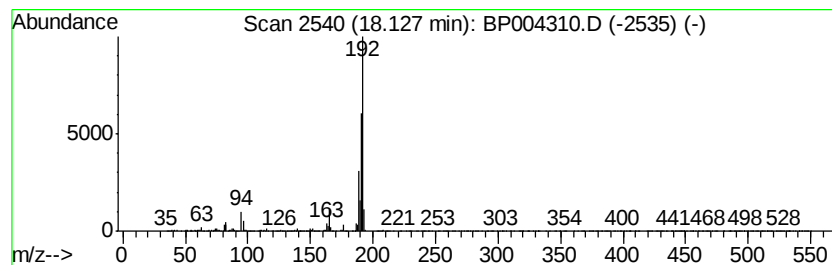
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Anthracene, 9-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.13	4.44 ng/ul	841381	Phenanthrene-d10	17.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
2		5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	87
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	86
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	83
5		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	81



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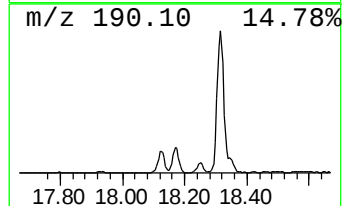
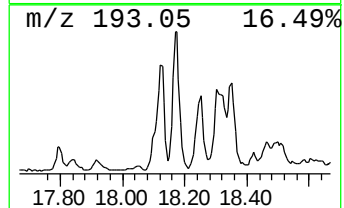
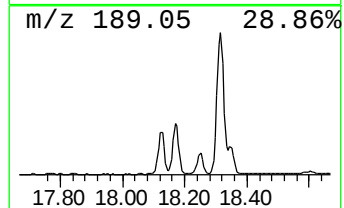
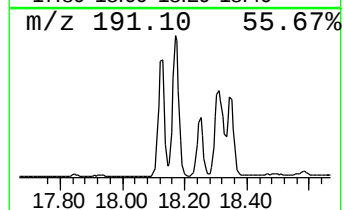
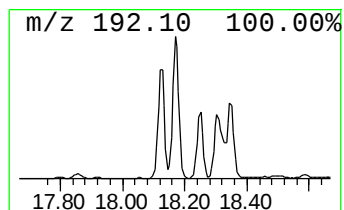
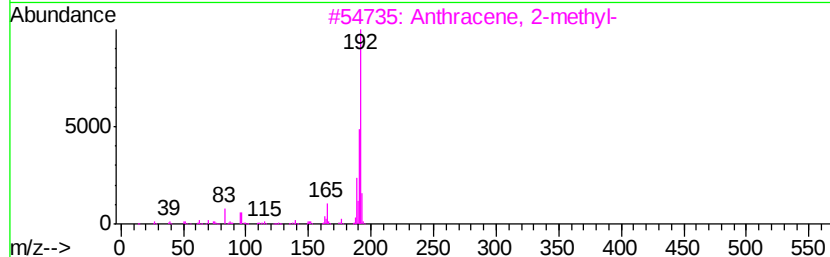
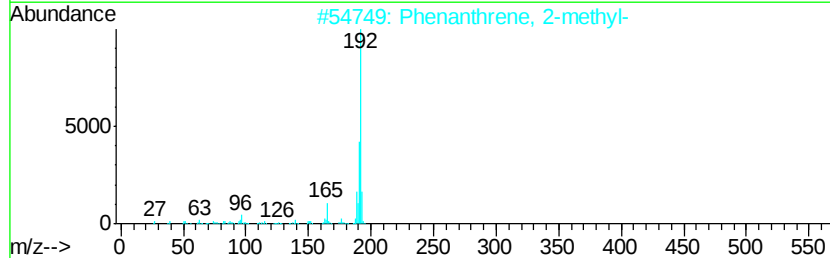
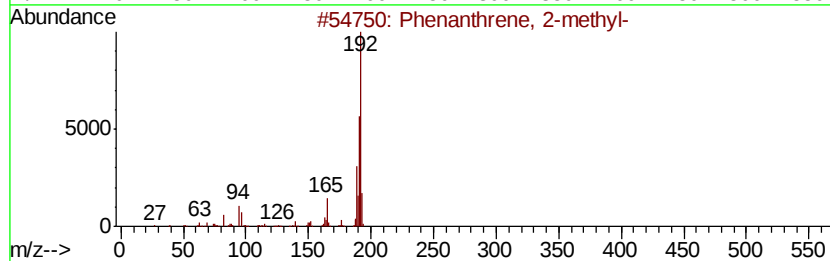
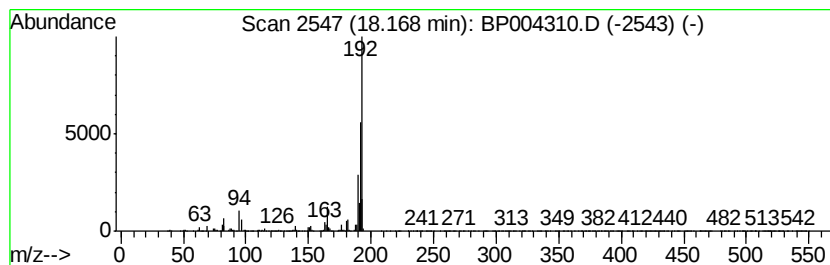
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.17	5.80 ng/ul	1098370	Phenanthrene-d10	17.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
3		Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4		Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
5		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95



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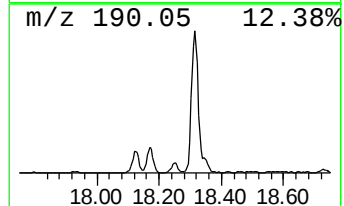
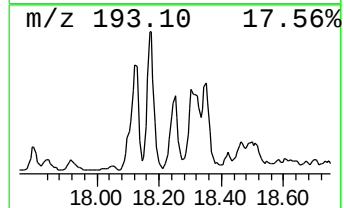
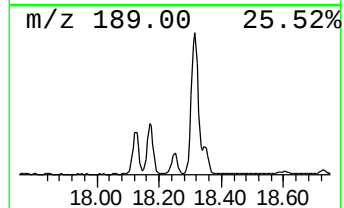
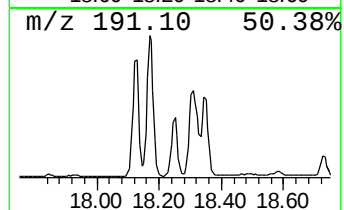
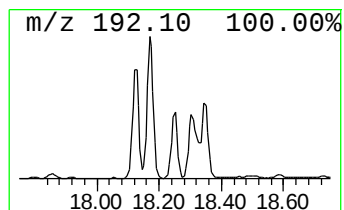
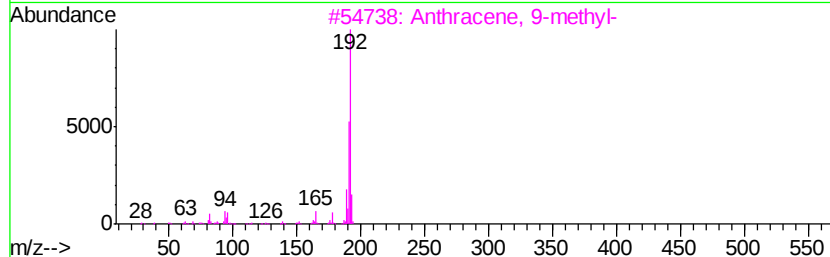
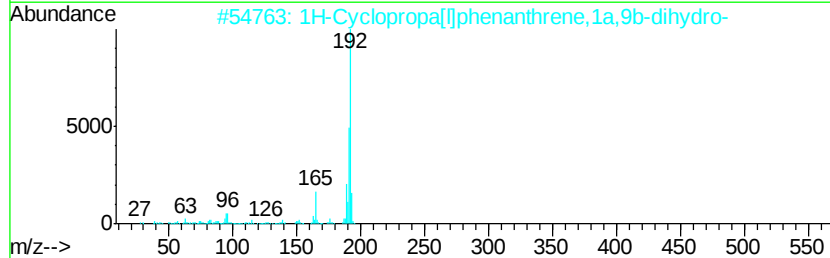
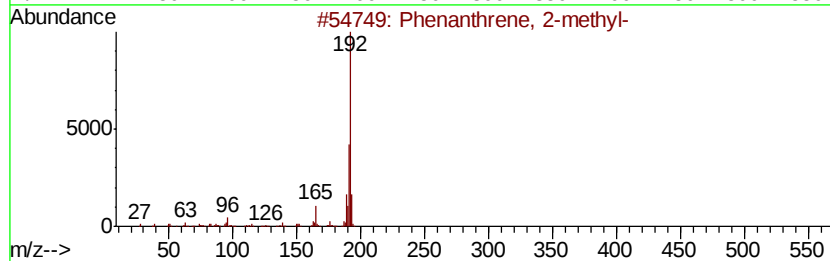
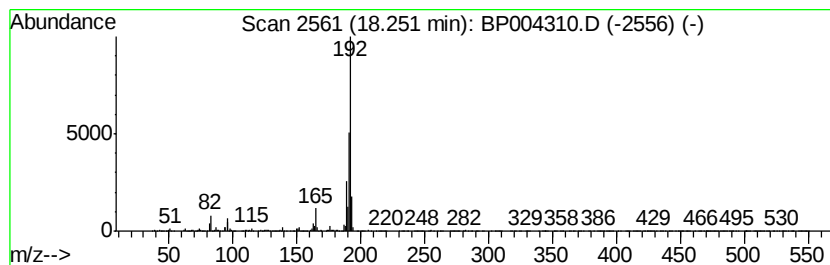
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1H-Cyclopropa[l]phenanthren... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.25	2.28 ng/ul	431444	Phenanthrene-d10	17.26	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	97
2	1H-Cyclopropa[ll]phenanthrene,1a,...	192	C15H12	000949-41-7	96
3	Anthracene, 9-methyl-	192	C15H12	000779-02-2	96
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
5	Anthracene, 2-methyl-	192	C15H12	000613-12-7	95



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\
Data File : BP004310.D
Acq On : 15 Dec 2020 18:34
Operator : CG/JU
Sample : L5001-13DL 10X
Misc :
ALS Vial : 45 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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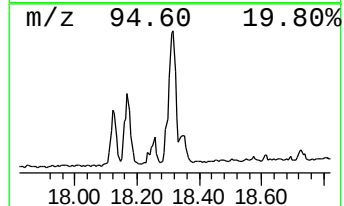
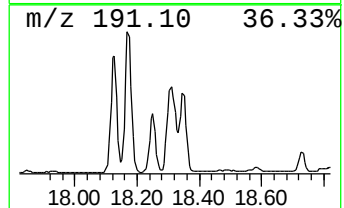
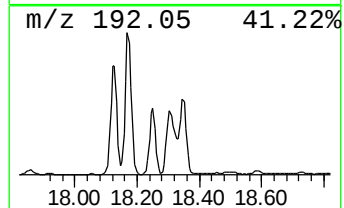
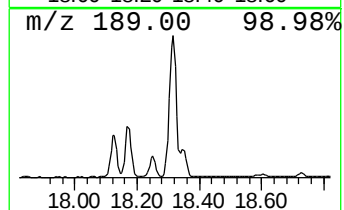
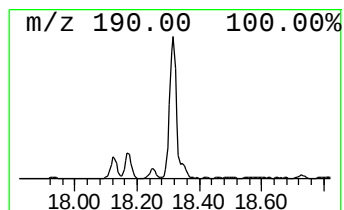
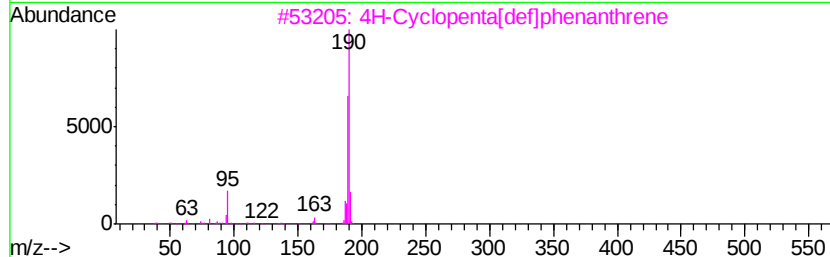
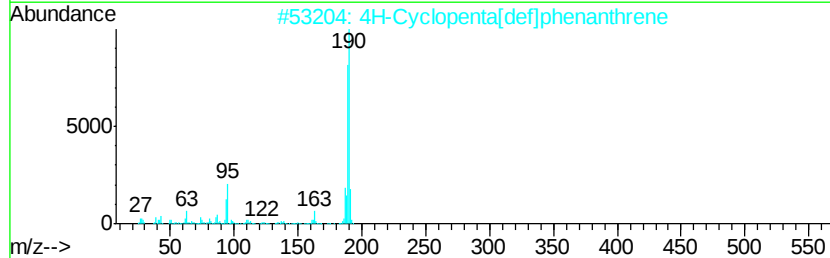
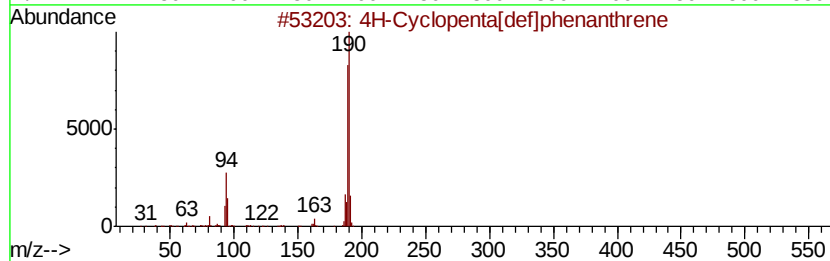
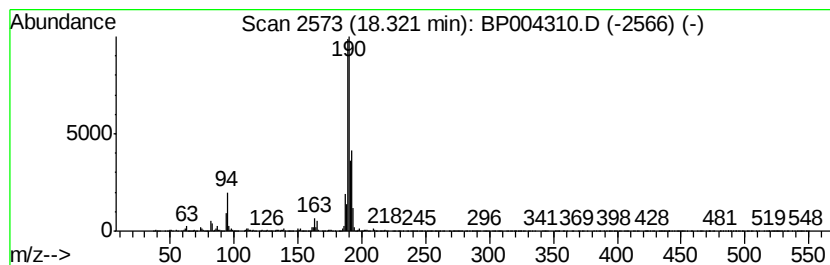
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 4H-Cyclopenta[def]phenanthrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.32	7.39 ng/ul	1399910	Phenanthrene-d10	17.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93
2		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
3		4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	62
4		6H-Cyclobuta[flk]phenanthrene	190	C15H10	083469-43-6	45
5		Methyl diselenide	190	C2H6Se2	007101-31-7	43



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\
Data File : BP004310.D
Acq On : 15 Dec 2020 18:34
Operator : CG/JU
Sample : L5001-13DL 10X
Misc :
ALS Vial : 45 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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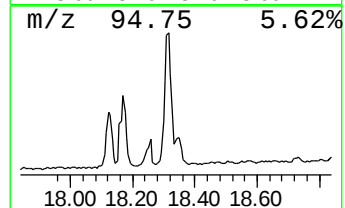
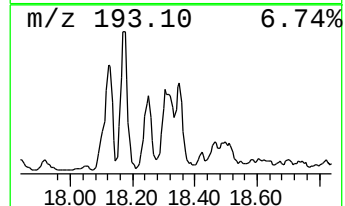
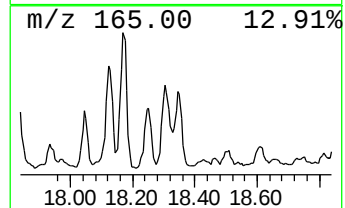
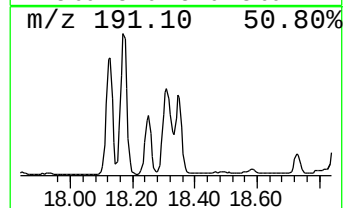
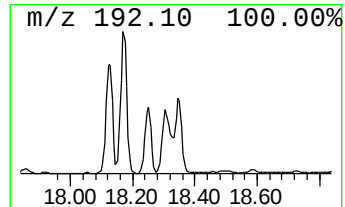
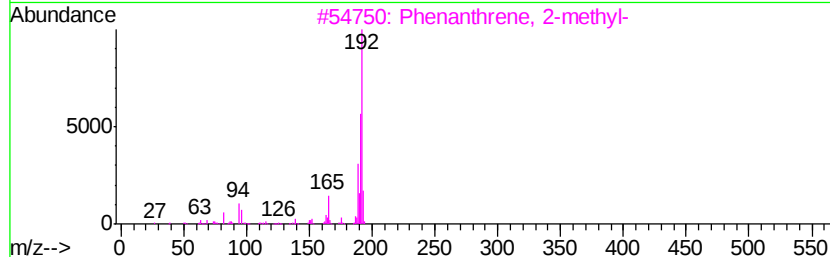
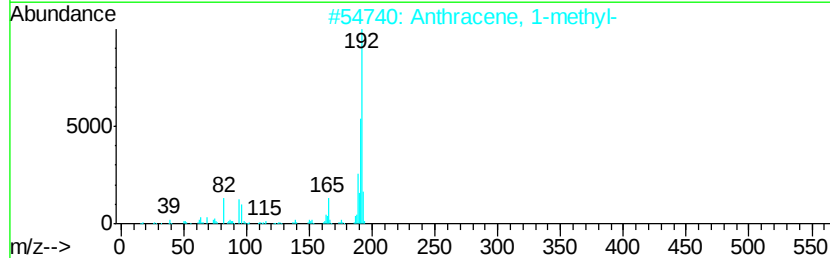
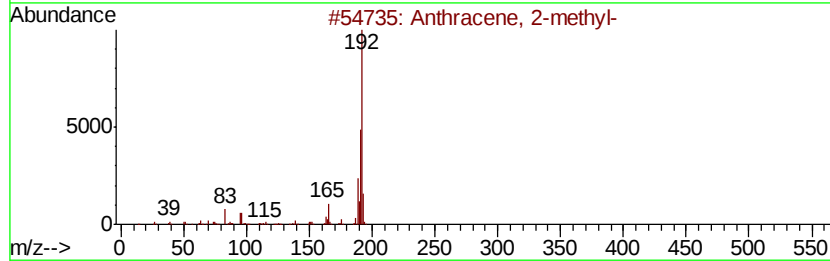
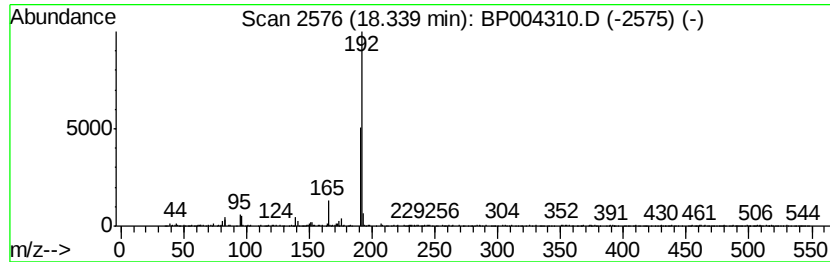
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.34	2.48 ng/ul	469549	Phenanthrene-d10	17.26

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
2		Anthracene, 1-methyl-	192	C15H12	000610-48-0	91
3		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	87
4		Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	87
5		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\
Data File : BP004310.D
Acq On : 15 Dec 2020 18:34
Operator : CG/JU
Sample : L5001-13DL 10X
Misc :
ALS Vial : 45 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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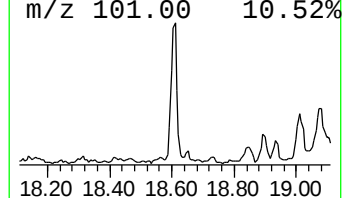
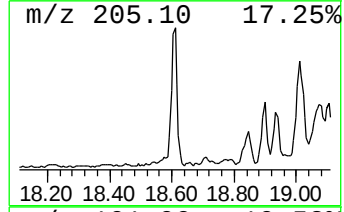
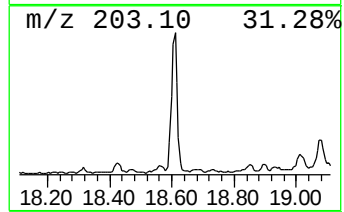
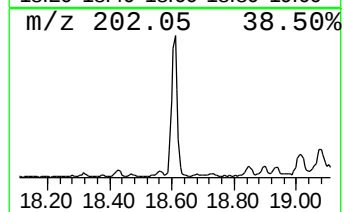
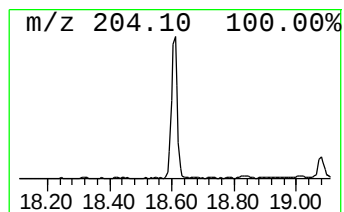
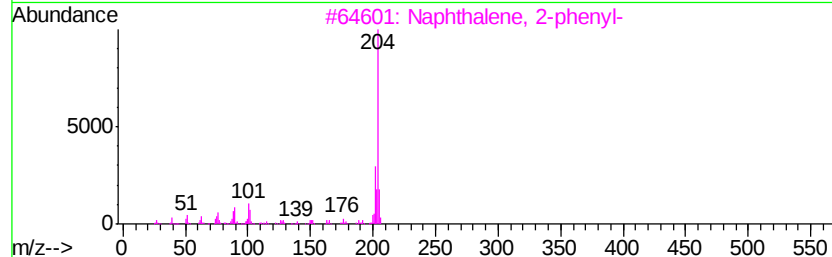
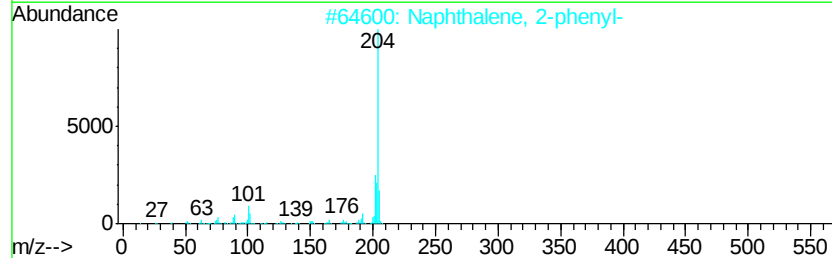
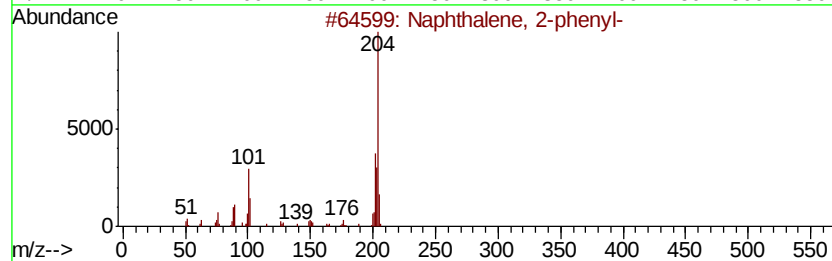
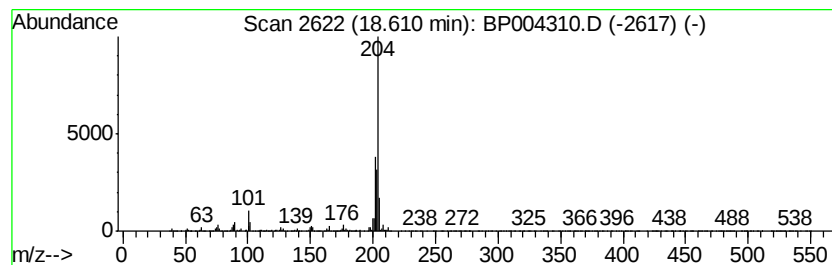
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Naphthalene, 2-phenyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.61	3.18 ng/ul	601758	Phenanthrene-d10	17.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	86
2		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	86
3		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	64
4		Tricyclo[8.2.2.2(4,7)]hexadeca-2...	204	C16H12	006572-60-7	62
5		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	60



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\
Data File : BP004310.D
Acq On : 15 Dec 2020 18:34
Operator : CG/JU
Sample : L5001-13DL 10X
Misc :
ALS Vial : 45 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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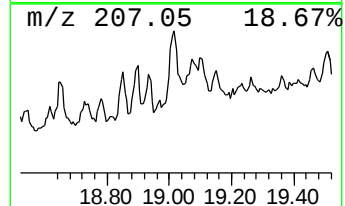
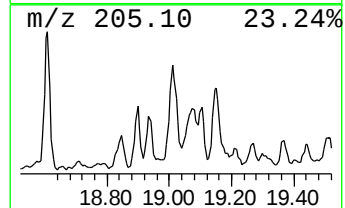
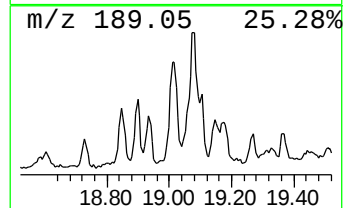
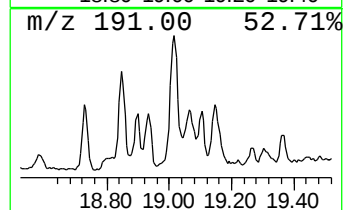
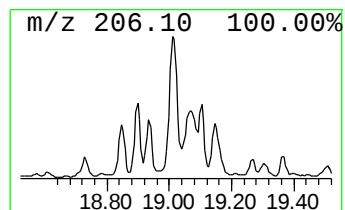
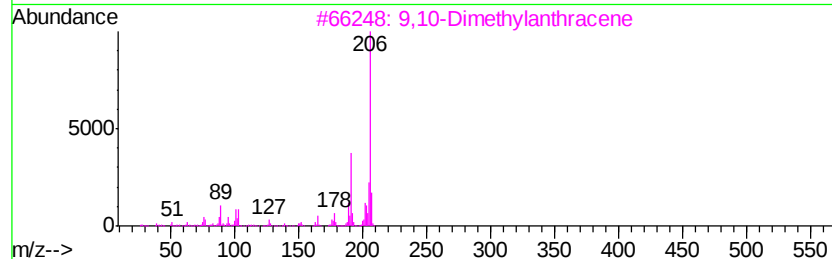
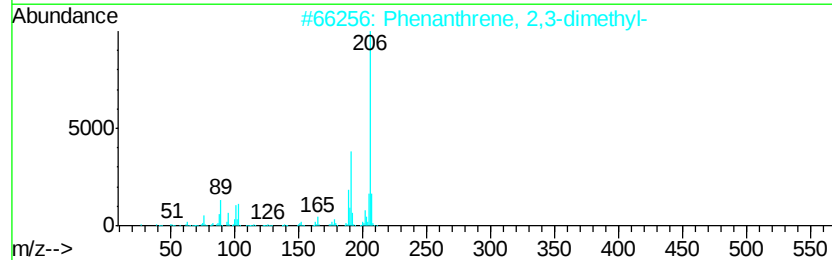
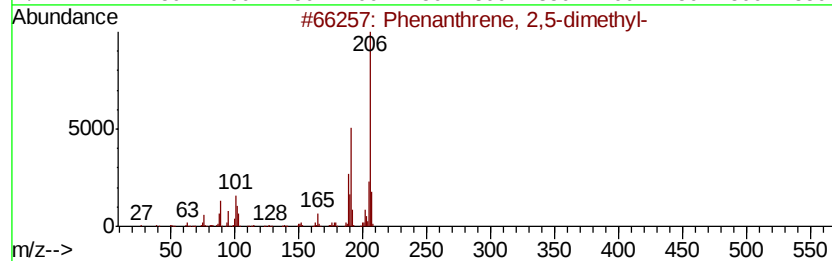
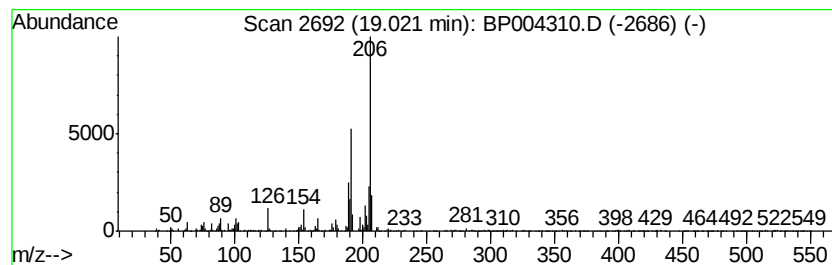
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Phenanthrene, 2,5-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.02	2.63 ng/ul	498080	Phenanthrene-d10	17.26

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	93
2		Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	87
3		9,10-Dimethylantracene	206	C16H14	000781-43-1	87
4		9,10-Dimethylantracene	206	C16H14	000781-43-1	87
5		Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\
Data File : BP004310.D
Acq On : 15 Dec 2020 18:34
Operator : CG/JU
Sample : L5001-13DL 10X
Misc :
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Instrument :
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ClientSampleId :
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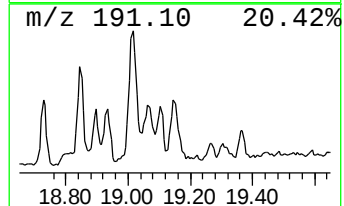
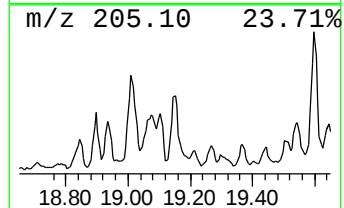
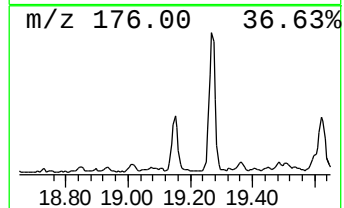
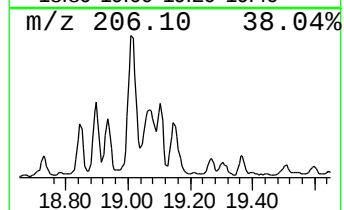
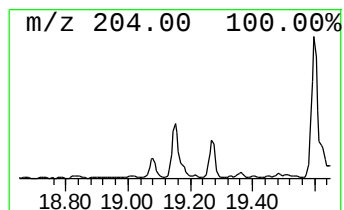
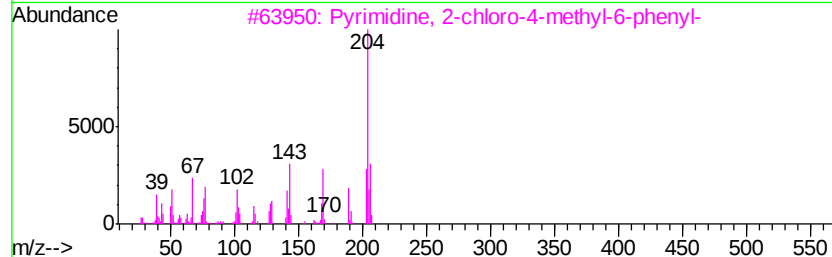
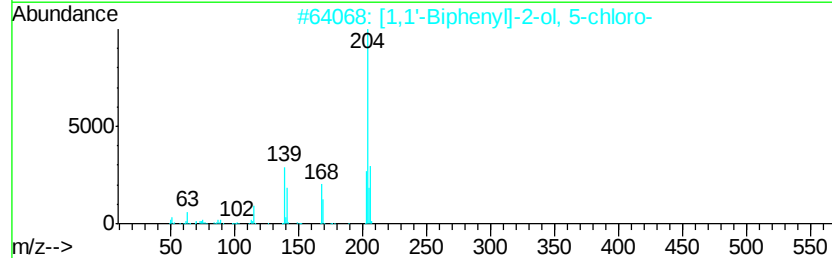
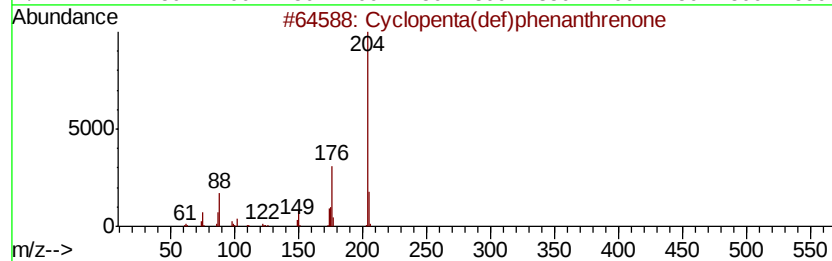
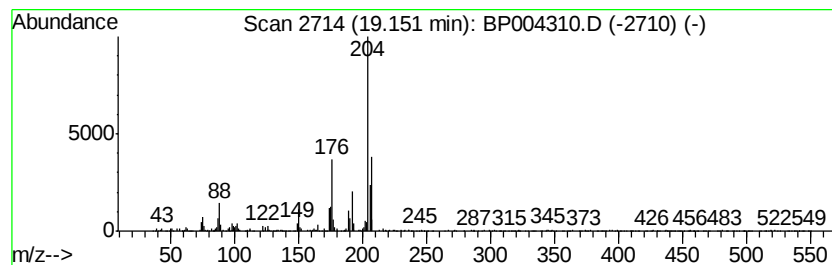
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Cyclopenta(def)phenanthrenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.15	2.25 ng/ul	425372	Phenanthrene-d10	17.26

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	92
2		[1,1'-Biphenyl]-2-ol, 5-chloro-	204	C12H9ClO	000607-12-5	50
3		Pvrimidine, 2-chloro-4-methyl-6-...	204	C11H9ClN2	032785-40-3	50
4		Benzene, 1-chloro-3-phenoxy-	204	C12H9ClO	006452-49-9	50
5		1,2,4,8-Tetramethylbicyclo[6.3.0...	204	C15H24	137235-51-9	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\
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Sample : L5001-13DL 10X
Misc :
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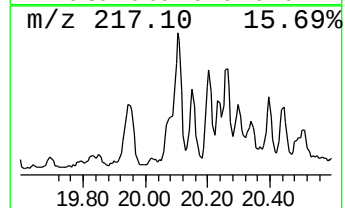
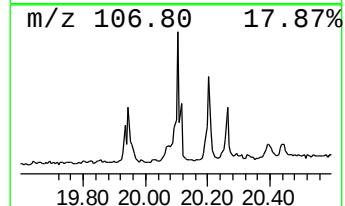
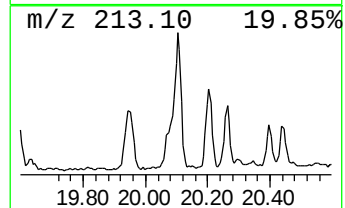
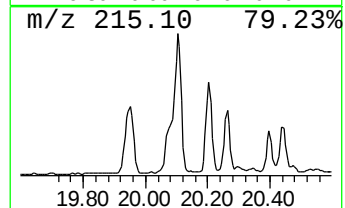
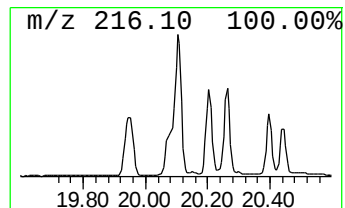
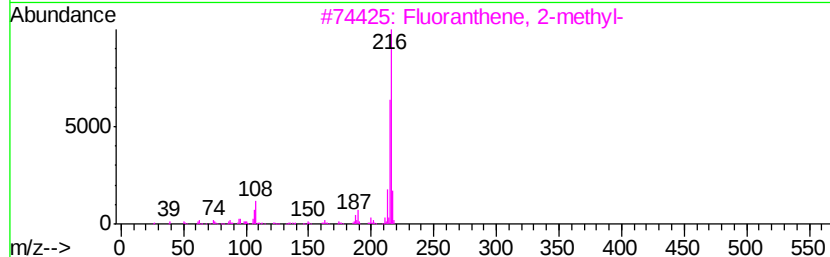
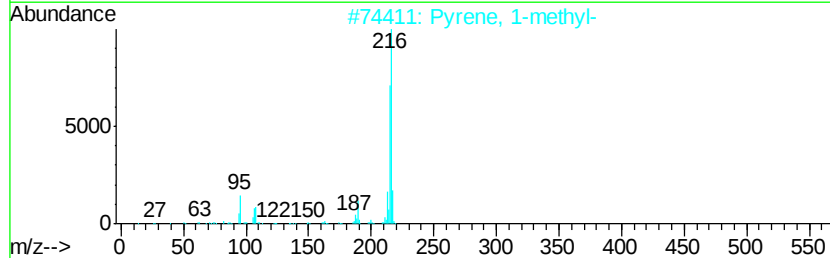
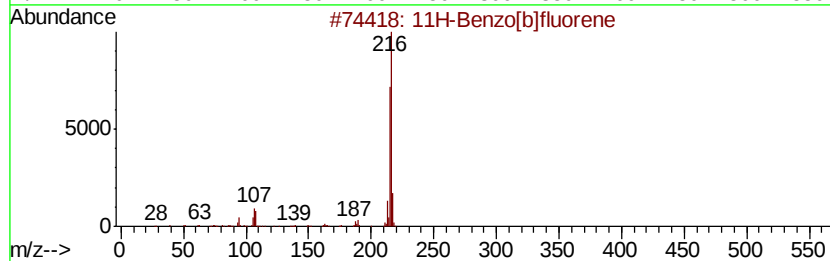
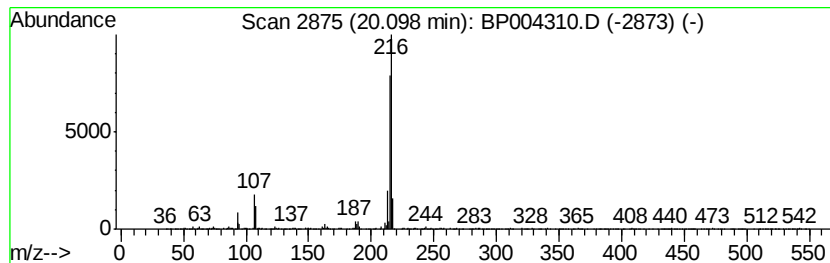
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 11H-Benzo[b]fluorene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.10	2.11 ng/ul	885249	Chrysene-d12	21.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		11H-Benzo[b]fluorene	216 C17H12	000243-17-4 93
2		Pyrene, 1-methyl-	216 C17H12	002381-21-7 81
3		Fluoranthene, 2-methyl-	216 C17H12	033543-31-6 76
4		11H-Benzo[a]fluorene	216 C17H12	000238-84-6 68
5		Fluoranthene, 2-methyl-	216 C17H12	033543-31-6 68



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121420\
Data File : BP004310.D
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Operator : CG/JU
Sample : L5001-13DL 10X
Misc :
ALS Vial : 45 Sample Multiplier: 1

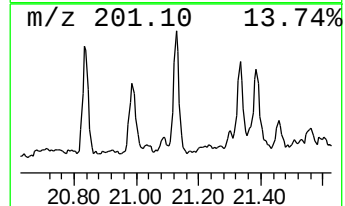
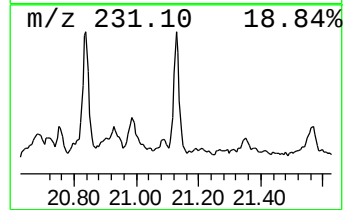
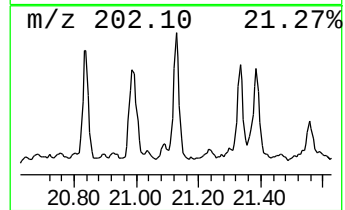
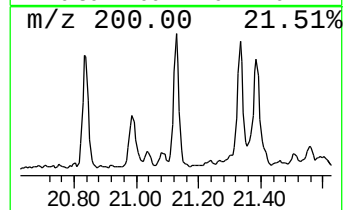
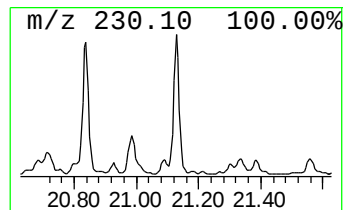
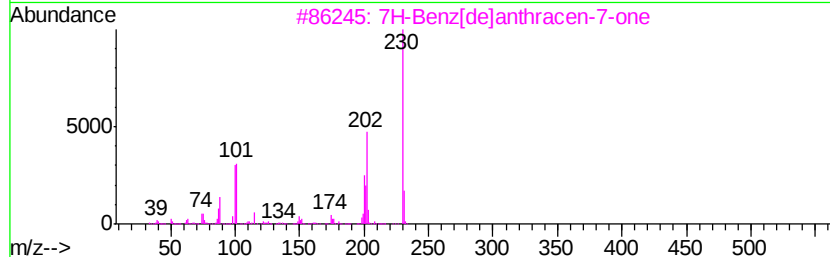
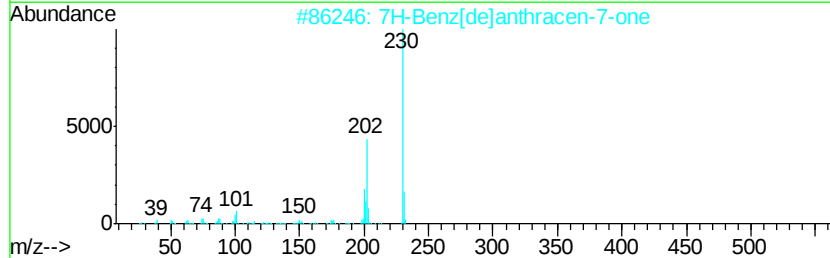
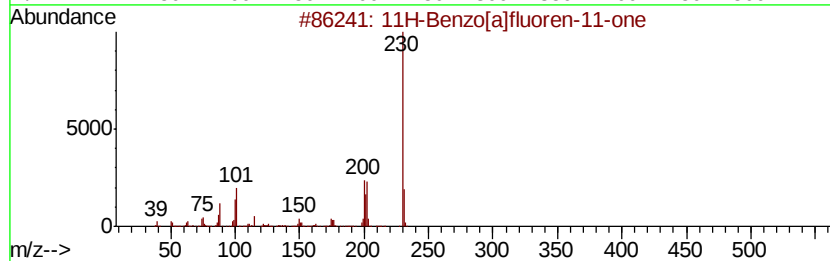
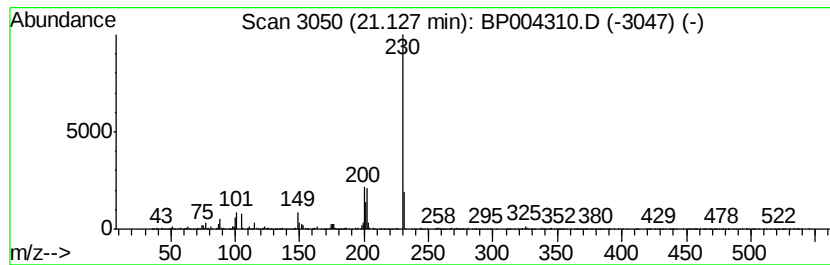
Instrument :
BNA_P
ClientSampleId :
A54D1DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 11H-Benzo[a]fluoren-11-one Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD		R.T.		
21.13	2.14 ng/ul	897858	Chrysene-d12		21.35		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	97
2			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	90
3			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	87
4			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	74
5			1-Pyrenecarboxaldehyde	230	C17H10O	003029-19-4	72



Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP121420\
 Data File : BP004310.D
 Acq On : 15 Dec 2020 18:34
 Operator : CG/JU
 Sample : L5001-13DL 10X
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A54D1DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP121020MA.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Anthracene, 9-met...	18.13	4.4	ng/ul	841381	4	17.26	3787780	20.0
Phenanthrene, 2-m...	18.17	5.8	ng/ul	1098370	4	17.26	3787780	20.0
1H-Cyclopropa[l]p...	18.25	2.3	ng/ul	431444	4	17.26	3787780	20.0
4H-Cyclopenta[def...	18.32	7.4	ng/ul	1399910	4	17.26	3787780	20.0
Anthracene, 2-met...	18.34	2.5	ng/ul	469549	4	17.26	3787780	20.0
Naphthalene, 2-ph...	18.61	3.2	ng/ul	601758	4	17.26	3787780	20.0
Phenanthrene, 2,5...	19.02	2.6	ng/ul	498080	4	17.26	3787780	20.0
Cyclopenta(def)ph...	19.15	2.3	ng/ul	425372	4	17.26	3787780	20.0
11H-Benzo[b]fluorene	20.10	2.1	ng/ul	885249	5	21.35	8390950	20.0
11H-Benzo[a]fluor...	21.13	2.1	ng/ul	897858	5	21.35	8390950	20.0