

Data Path : Z:\HPCHEM1\BNA G\DATA\BG051616\
 Data File : BG021916.D
 Acq On : 16 May 2016 21:25
 Operator : UM/SJ
 Sample : H3095-07
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 D9XE4

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:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG050916.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	3.229	62	66	74	rVB3	10585	18726	1.27%	0.107%
2	3.329	78	83	87	rBV6	3764	7014	0.48%	0.040%
3	3.546	115	120	125	rBV3	7062	13823	0.94%	0.079%
4	4.081	206	211	219	rBV	8302	15968	1.08%	0.091%
5	4.304	244	249	256	rBV2	3407	6104	0.41%	0.035%
6	5.332	418	424	429	rBV4	3542	6438	0.44%	0.037%
7	5.785	495	501	508	rBV5	3811	9527	0.65%	0.054%
8	6.155	560	564	570	rVV3	3717	7184	0.49%	0.041%
9	7.248	743	750	754	rBV3	2582	4760	0.32%	0.027%
10	7.318	754	762	775	rBV	112475	240136	16.30%	1.373%
11	7.489	784	791	800	rBV	100712	189589	12.87%	1.084%
12	7.700	817	827	836	rBV	134359	267902	18.19%	1.532%
13	7.776	836	840	843	rVV3	3656	7436	0.50%	0.043%
14	7.806	843	845	851	rVB3	5541	7750	0.53%	0.044%
15	8.170	893	907	921	rBV	156736	317702	21.57%	1.817%
16	8.276	921	925	932	rVB5	3015	5242	0.36%	0.030%
17	8.717	998	1000	1008	rVB5	2939	6539	0.44%	0.037%
18	8.811	1011	1016	1019	rBV3	2947	5559	0.38%	0.032%
19	8.875	1019	1027	1041	rBV	156806	325614	22.10%	1.862%
20	9.010	1045	1050	1055	rVB6	3829	7614	0.52%	0.044%
21	9.281	1087	1096	1098	rBV5	3700	8021	0.54%	0.046%
22	9.339	1098	1106	1113	rBV	109229	226094	15.35%	1.293%
23	9.486	1126	1131	1140	rVB6	3914	8282	0.56%	0.047%
24	10.062	1221	1229	1240	rBV	87239	190454	12.93%	1.089%
25	10.385	1278	1284	1289	rBV3	4046	7653	0.52%	0.044%
26	10.608	1314	1322	1333	rVV	166577	347707	23.60%	1.988%
27	10.932	1374	1377	1379	rBV2	4821	6749	0.46%	0.039%
28	10.990	1379	1387	1394	rBV	255977	512159	34.77%	2.928%
29	11.125	1402	1410	1423	rBV	82858	197918	13.44%	1.132%
30	11.983	1550	1556	1564	rVV3	10895	19985	1.36%	0.114%
31	12.371	1615	1622	1629	rBV5	10354	21310	1.45%	0.122%
32	12.565	1649	1655	1658	rVV2	3757	7230	0.49%	0.041%
33	12.635	1660	1667	1674	rVB	18858	34391	2.33%	0.197%
34	12.847	1696	1703	1710	rBV	16770	33442	2.27%	0.191%

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35	13.317	1777	1783	1787	rBV3	8570	13347	0.91%	0.076%
36	13.393	1791	1796	1801	rBV5	5008	9159	0.62%	0.052%
37	13.593	1825	1830	1834	rBV3	13995	24588	1.67%	0.141%
38	13.675	1840	1844	1849	rBV6	3468	7132	0.48%	0.041%
39	13.969	1887	1894	1901	rVB2	8744	21443	1.46%	0.123%
40	14.116	1913	1919	1924	rBV3	30308	51576	3.50%	0.295%
41	14.187	1924	1931	1936	rBV	318275	562239	38.17%	3.215%
42	14.234	1936	1939	1947	rVB	91572	159454	10.82%	0.912%
43	14.345	1955	1958	1962	rBV5	7703	13630	0.93%	0.078%
44	14.492	1976	1983	1997	rBV	450419	842467	57.19%	4.817%
45	14.598	1997	2001	2003	rBV3	6681	9407	0.64%	0.054%
46	14.663	2008	2012	2019	rVB2	24811	36268	2.46%	0.207%
47	14.739	2019	2025	2028	rBV3	23659	50716	3.44%	0.290%
48	14.792	2028	2034	2042	rVV2	412413	751485	51.01%	4.297%
49	14.856	2042	2045	2052	rVB5	16264	30161	2.05%	0.172%
50	14.992	2058	2068	2074	rBV2	65879	189932	12.89%	1.086%
51	15.039	2074	2076	2085	rVV10	16698	34162	2.32%	0.195%
52	15.185	2095	2101	2108	rVB2	17585	34719	2.36%	0.199%
53	15.262	2110	2114	2122	rVB8	8804	22361	1.52%	0.128%
54	15.403	2133	2138	2143	rBV5	11060	20277	1.38%	0.116%
55	15.450	2143	2146	2151	rVV4	6040	9869	0.67%	0.056%
56	15.573	2161	2167	2170	rBV6	15549	29258	1.99%	0.167%
57	15.608	2170	2173	2178	rVB	41250	59577	4.04%	0.341%
58	15.667	2178	2183	2184	rBV3	6279	7853	0.53%	0.045%
59	15.785	2193	2203	2209	rBV	594601	1004289	68.17%	5.742%
60	15.826	2209	2210	2217	rVV5	24470	32662	2.22%	0.187%
61	15.902	2217	2223	2232	rVB	79704	126533	8.59%	0.723%
62	16.020	2237	2243	2247	rVB6	19300	34569	2.35%	0.198%
63	16.125	2254	2261	2265	rBV4	12053	26733	1.81%	0.153%
64	16.237	2277	2280	2282	rBV4	6154	7130	0.48%	0.041%
65	16.272	2284	2286	2296	rVB3	10732	20442	1.39%	0.117%
66	16.472	2315	2320	2322	rBV	48627	72109	4.89%	0.412%
67	16.619	2341	2345	2347	rBV5	5225	8280	0.56%	0.047%
68	16.813	2376	2378	2380	rBV3	5162	5065	0.34%	0.029%
69	16.983	2403	2407	2413	rVB8	8555	15285	1.04%	0.087%
70	17.071	2415	2422	2424	rBV5	13524	32141	2.18%	0.184%
71	17.195	2438	2443	2448	rBV	20483	34870	2.37%	0.199%

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72	17.265	2450	2455	2461	rVV	42330	68524	4.65%	0.392%
73	17.318	2461	2464	2468	rVB4	10687	13825	0.94%	0.079%
74	17.353	2468	2470	2475	rVB3	12133	14929	1.01%	0.085%
75	17.406	2475	2479	2486	rVB8	16660	24757	1.68%	0.142%
76	17.471	2487	2490	2492	rBV4	7696	7878	0.53%	0.045%
77	17.541	2495	2502	2506	rBV2	488158	914981	62.11%	5.232%
78	17.583	2506	2509	2513	rVV	223603	335786	22.79%	1.920%
79	17.635	2513	2518	2530	rVB2	548068	976689	66.30%	5.584%
80	17.941	2564	2570	2576	rBV4	21209	53401	3.62%	0.305%
81	18.006	2578	2581	2585	rVB2	29451	38919	2.64%	0.223%
82	18.117	2598	2600	2607	rVB6	14380	21995	1.49%	0.126%
83	18.346	2633	2639	2644	rBV9	19554	44554	3.02%	0.255%
84	18.405	2644	2649	2654	rVV2	90015	159728	10.84%	0.913%
85	18.458	2654	2658	2666	rVB2	67055	138031	9.37%	0.789%
86	18.599	2677	2682	2687	rBV2	61871	130989	8.89%	0.749%
87	18.693	2694	2698	2705	rBV4	32141	57781	3.92%	0.330%
88	18.905	2730	2734	2737	rBV	37514	52897	3.59%	0.302%
89	18.946	2738	2741	2746	rVB3	19839	31420	2.13%	0.180%
90	19.316	2800	2804	2809	rBV	68543	96117	6.52%	0.550%
91	19.463	2824	2829	2835	rBV2	35812	69108	4.69%	0.395%
92	19.580	2844	2849	2855	rBV	468372	709320	48.15%	4.056%
93	19.915	2899	2906	2909	rBV	714585	1267542	86.04%	7.247%
94	20.420	2989	2992	2998	rVB2	91811	143838	9.76%	0.822%
95	21.431	3161	3164	3170	rVB2	113740	152705	10.37%	0.873%
96	21.690	3202	3208	3216	rVB	929916	1473137	100.00%	8.423%
97	21.825	3223	3231	3236	rBV2	565555	1340305	90.98%	7.663%
98	24.104	3613	3619	3626	rBV	137399	430557	29.23%	2.462%
99	24.939	3755	3761	3768	rBV2	204354	485198	32.94%	2.774%
100	25.168	3793	3800	3811	rVB2	250647	731539	49.66%	4.183%

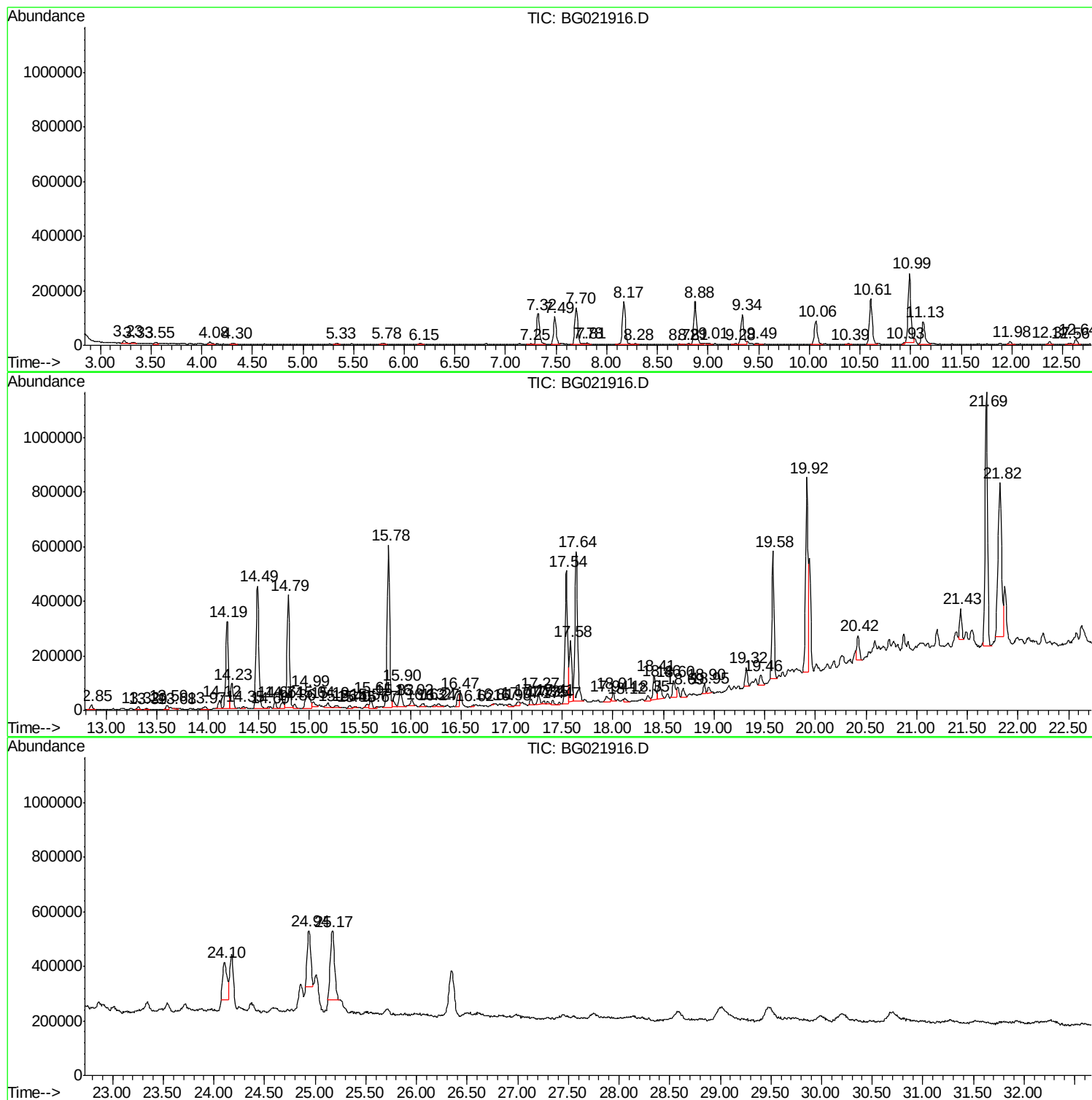
Sum of corrected areas: 17489660

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG050916.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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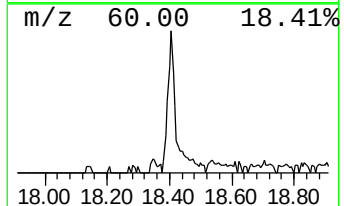
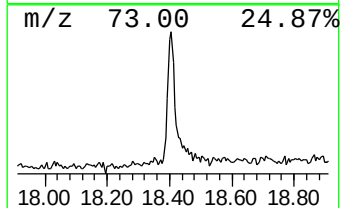
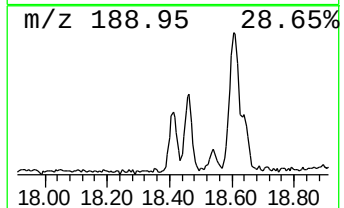
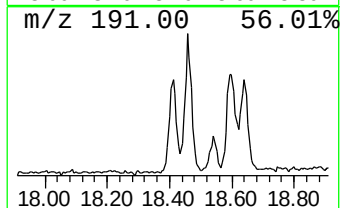
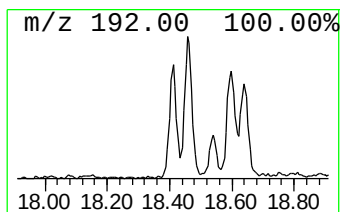
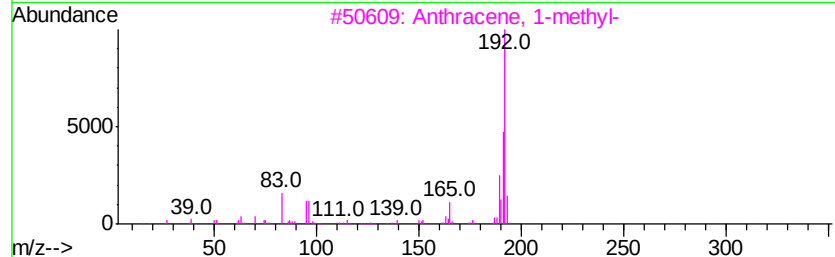
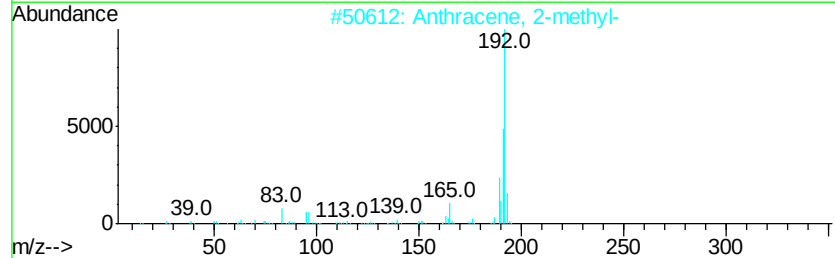
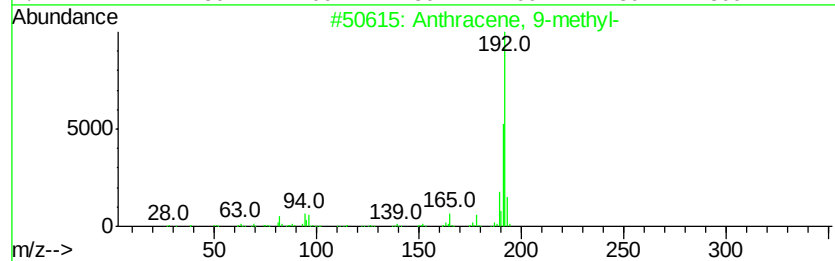
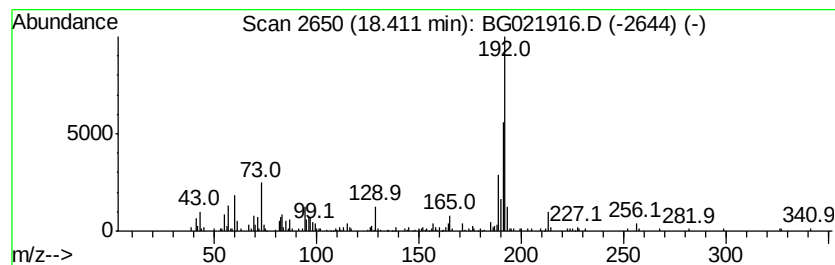
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG050916.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.41	3.49 ng/ul	159728	Phenanthrene-d10	17.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 9-methyl-	192	C15H12	000779-02-2	92
2		Anthracene, 2-methyl-	192	C15H12	000613-12-7	87
3		Anthracene, 1-methyl-	192	C15H12	000610-48-0	83
4		Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	83
5		1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	83



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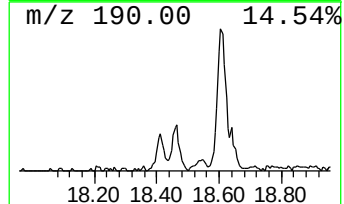
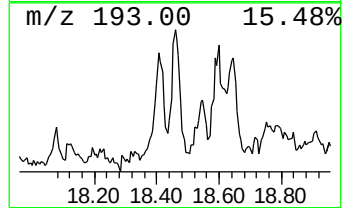
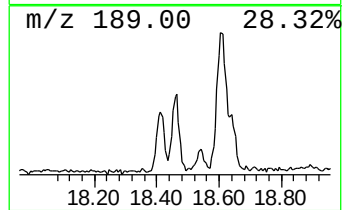
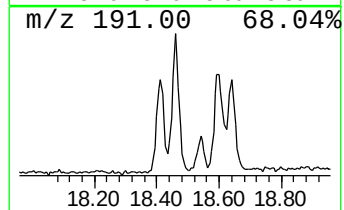
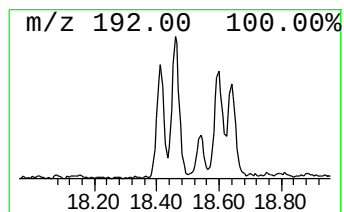
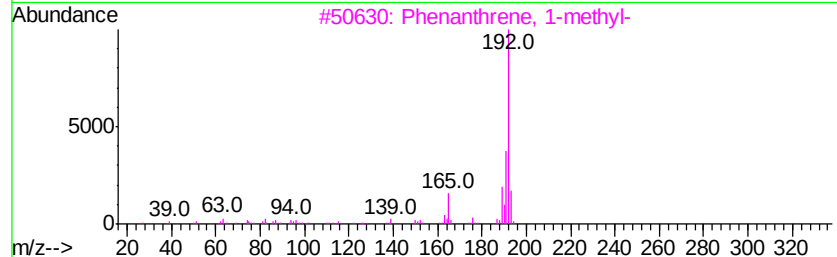
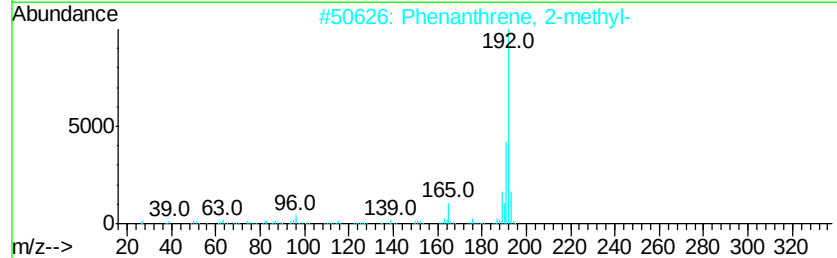
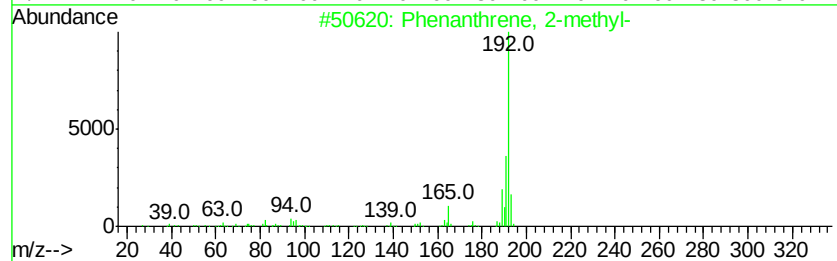
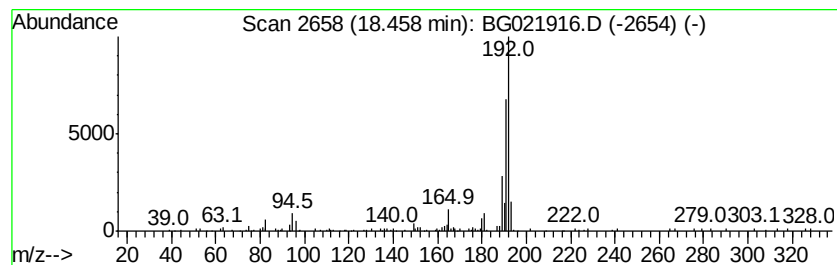
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG050916.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.46	3.02 ng/ul	138031	Phenanthrene-d10	17.54	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	86
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	83
3	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	83
4	1H-Cyclopropa[1,1]phenanthrene,1a,...	192	C15H12	000949-41-7	83
5	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	81



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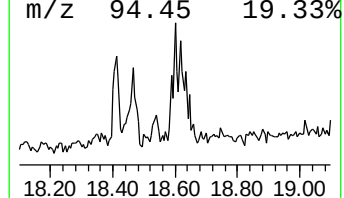
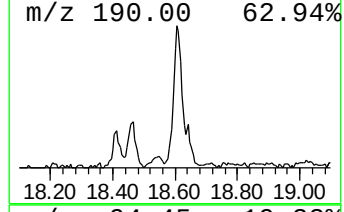
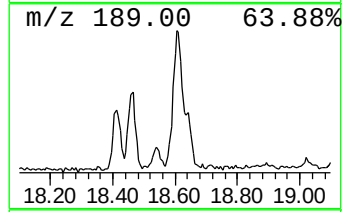
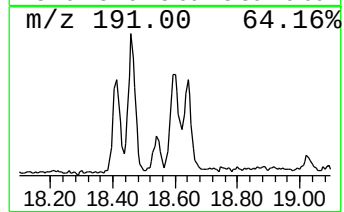
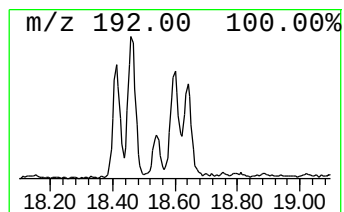
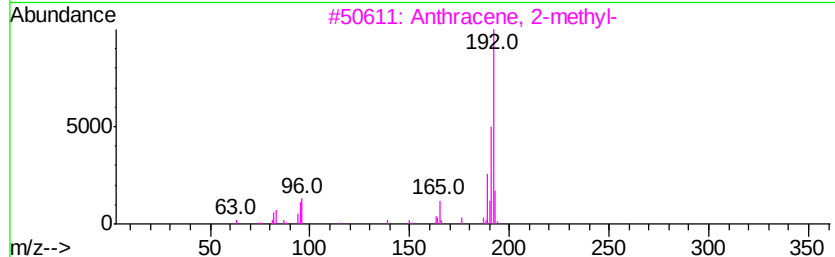
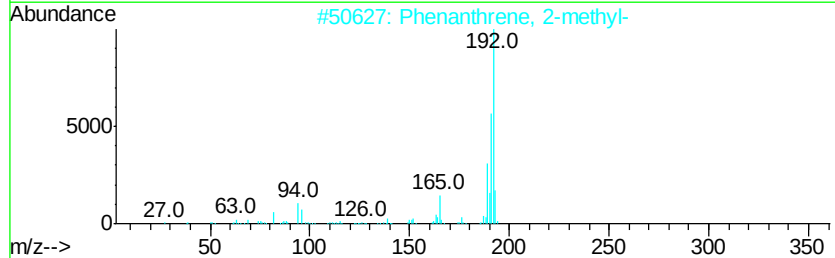
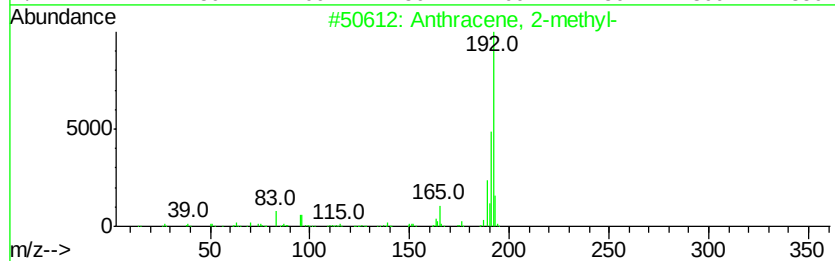
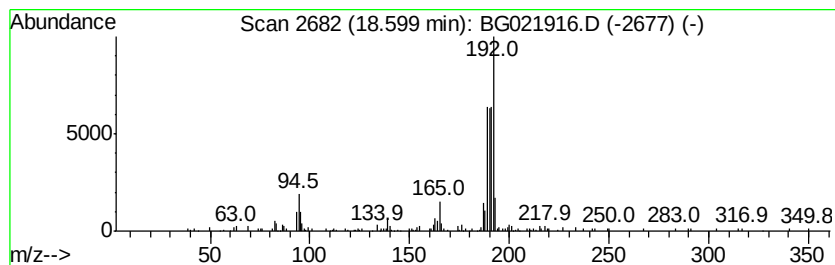
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.60	2.86 ng/ul	130989	Phenanthrene-d10	17.54

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 2-methyl-	192	C15H12	000613-12-7	60
2			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	60
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	60
4			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	55
5			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	55



Data Path : Z:\HPCHEM1\BNA G\DATA\BG051616\
Data File : BG021916.D
Acq On : 16 May 2016 21:25
Operator : UM/SJ
Sample : H3095-07
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
D9XE4

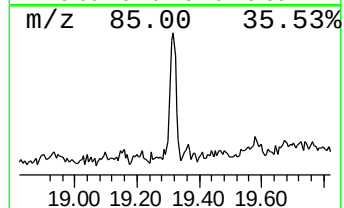
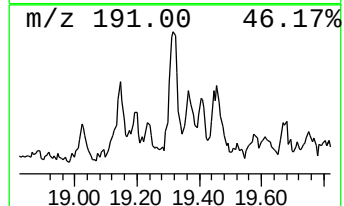
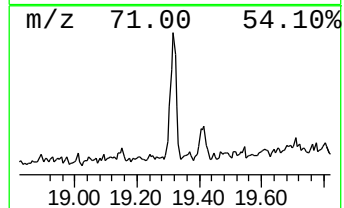
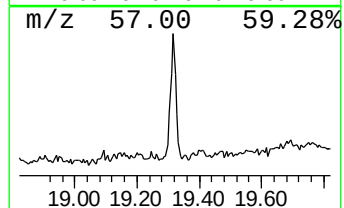
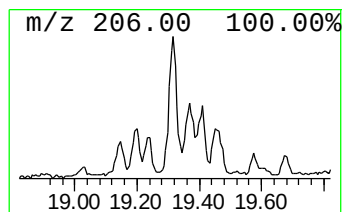
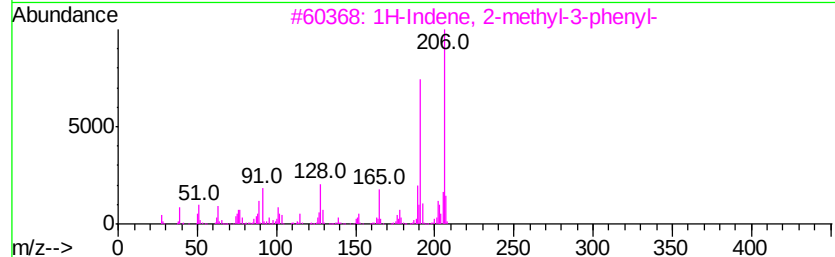
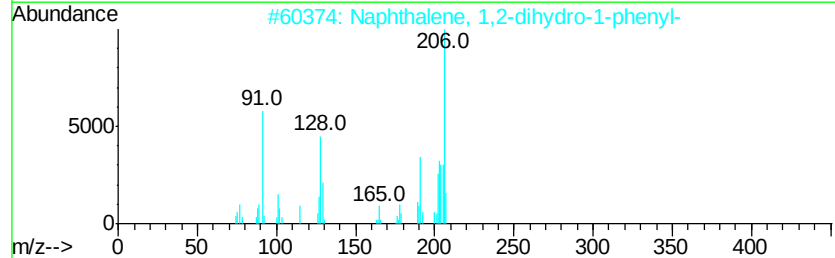
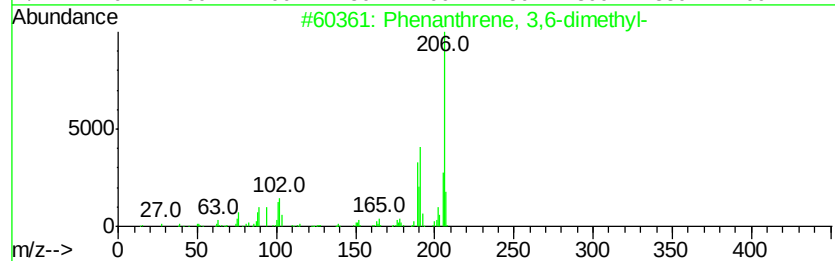
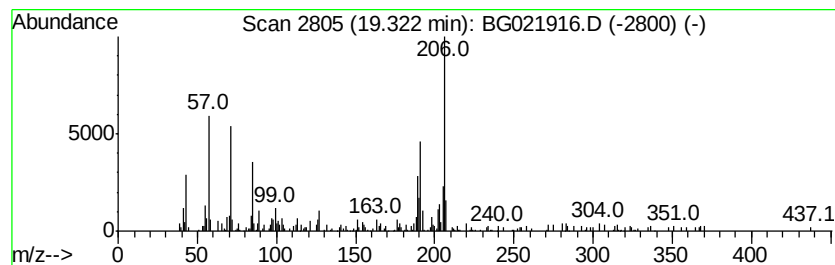
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG050916.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Phenanthrene, 3,6-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.32	2.10 ng/ul	96117	Phenanthrene-d10	17.54

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	83
2		Naphthalene, 1,2-dihydro-1-phenyl-	206	C16H14	016606-46-5	74
3		1H-Indene, 2-methyl-3-phenyl-	206	C16H14	035099-60-6	70
4		di-p-Tolylacetylene	206	C16H14	002789-88-0	64
5		Naphthalene, 1,2-dihydro-4-phenyl-	206	C16H14	007469-40-1	64



Data Path : Z:\HPCHEM1\BNA G\DATA\BG051616\
Data File : BG021916.D
Acq On : 16 May 2016 21:25
Operator : UM/SJ
Sample : H3095-07
Misc :
ALS Vial : 16 Sample Multiplier: 1

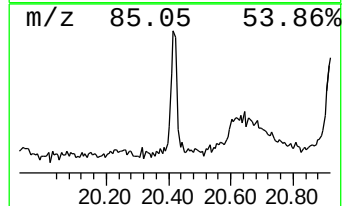
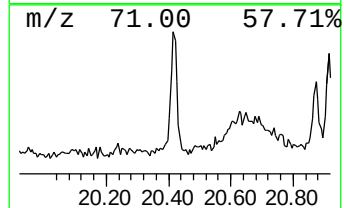
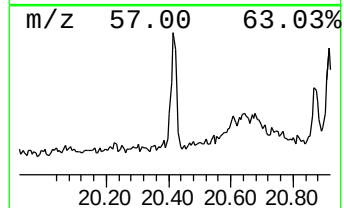
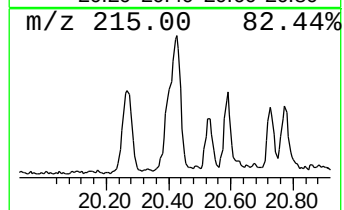
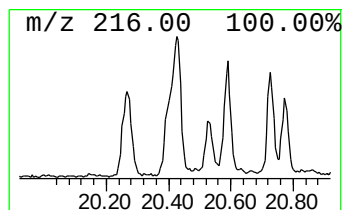
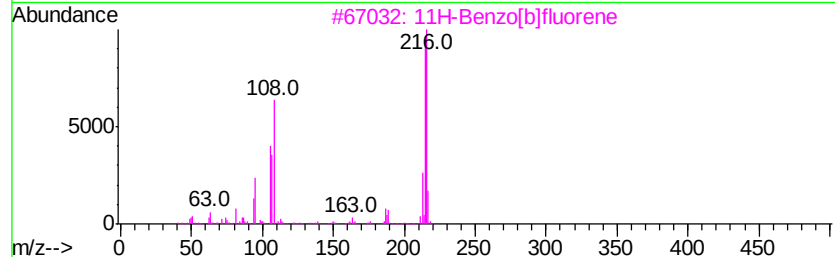
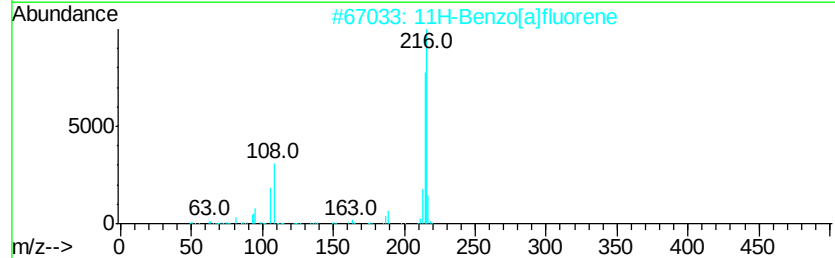
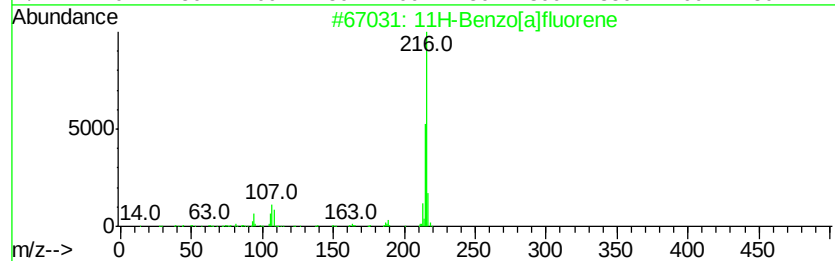
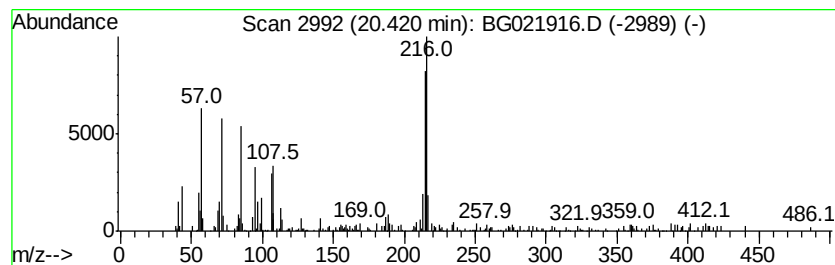
Instrument :
BNA_G
ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG050916.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 11H-Benzo[a]fluorene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.42	2.15 ng/ul	143838	Chrysene-d12	21.82	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	74
2	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	60
3	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	58
4	1,3,5-Triazine-2,4,6-triamine, N...	216	C10H12N6	046731-79-7	58
5	Pyrene, 1-methyl-	216	C17H12	002381-21-7	53



Data Path : Z:\HPCHEM1\BNA G\DATA\BG051616\
Data File : BG021916.D
Acq On : 16 May 2016 21:25
Operator : UM/SJ
Sample : H3095-07
Misc :
ALS Vial : 16 Sample Multiplier: 1

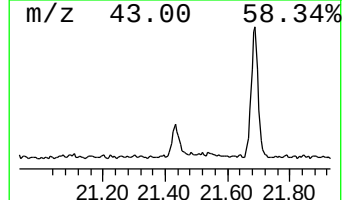
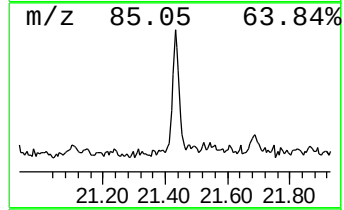
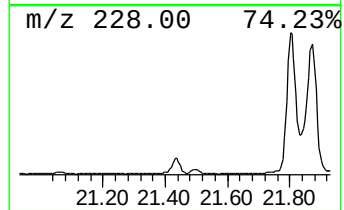
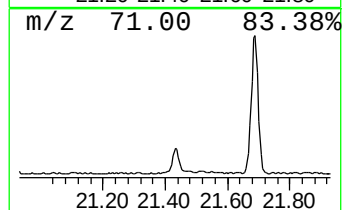
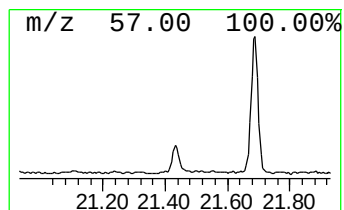
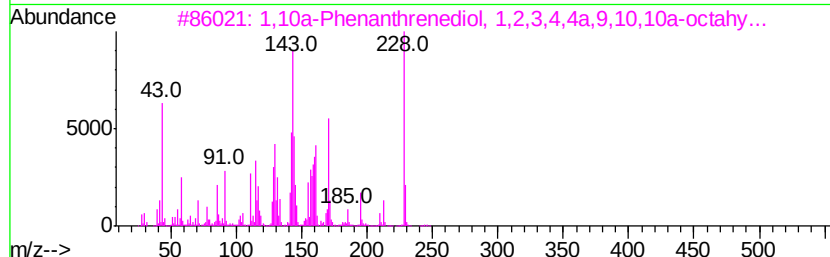
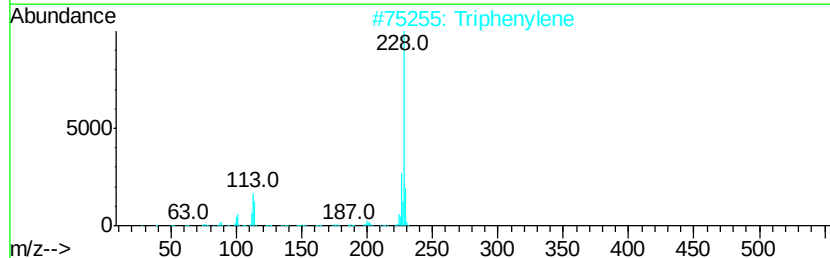
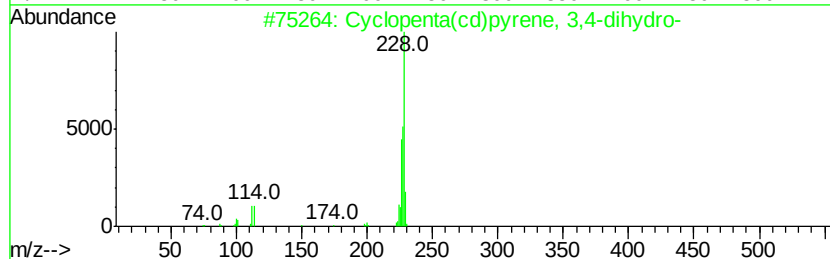
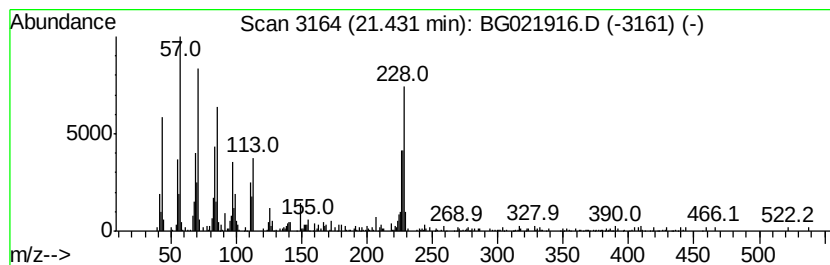
Instrument :
BNA_G
ClientSampleId :
D9XE4

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG050916.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.43	2.28 ng/ul	152705	Chrysene-d12	21.82
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclopenta(cd)pyrene, 3,4-dihydro-	228 C18H12	025732-74-5 38
2		Triphenylene	228 C18H12	000217-59-4 14
3		1,10a-Phenanthrenediol, 1,2,3,4,...	246 C16H22O2	1000197-42-1 14
4		5-Ethyl-5-acetyl-amino-6-imino-1,...	228 C8H12N4O2S	126552-74-7 14
5		2-Formylphenoxathin	228 C13H8O2S	037947-92-5 11



Data Path : Z:\HPCHEM1\BNA_G\DATA\BG051616\
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Acq On : 16 May 2016 21:25
Operator : UM/SJ
Sample : H3095-07
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
D9XE4

Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG050916.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Anthracene, 9-met...	18.41	3.5	ng/ul	159728	4	17.54	914981	20.0
Phenanthrene, 2-m...	18.46	3.0	ng/ul	138031	4	17.54	914981	20.0
Anthracene, 2-met...	18.60	2.9	ng/ul	130989	4	17.54	914981	20.0
Phenanthrene, 3,6...	19.32	2.1	ng/ul	96117	4	17.54	914981	20.0
11H-Benzo[a]fluorene	20.42	2.1	ng/ul	143838	5	21.82	1340310	20.0
unknown-01	21.43	2.3	ng/ul	152705	5	21.82	1340310	20.0