

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN070518\
 Data File : BN001974.D
 Acq On : 05 Jul 2018 19:06
 Operator : JU/SJ
 Sample : J3721-14DL 25X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 F1H00DL

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_N\METHODS\SOM-EPA-BN061918.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	3.022	3	6	13	rVB	194128	244867	1.05%	0.105%
2	6.881	655	662	672	rVB	81158	163392	0.70%	0.070%
3	7.052	685	691	698	rBV	66077	107929	0.46%	0.046%
4	7.246	717	724	731	rBV	80480	138019	0.59%	0.059%
5	7.710	796	803	811	rBV	2196422	3497265	14.94%	1.498%
6	8.410	917	922	929	rBV	92069	152265	0.65%	0.065%
7	8.857	992	998	1008	rBV	74377	134761	0.58%	0.058%
8	9.569	1114	1119	1125	rBV	53346	87677	0.37%	0.038%
9	10.104	1205	1210	1216	rBV2	92518	157746	0.67%	0.068%
10	10.481	1266	1274	1281	rBV	3276038	5499207	23.50%	2.355%
11	10.622	1290	1298	1303	rBV2	39497	71710	0.31%	0.031%
12	13.751	1821	1830	1834	rBV	193766	287376	1.23%	0.123%
13	13.792	1834	1837	1842	rVB2	95525	131860	0.56%	0.056%
14	14.028	1871	1877	1878	rBV	222100	287084	1.23%	0.123%
15	14.057	1878	1882	1894	rVB	509306	832172	3.56%	0.356%
16	14.340	1920	1930	1937	rBV2	4677815	7359183	31.44%	3.152%
17	15.334	2093	2099	2104	rBV	328005	482535	2.06%	0.207%
18	15.387	2104	2108	2115	rVB	61249	96349	0.41%	0.041%
19	15.610	2138	2146	2152	rBV8	27407	83380	0.36%	0.036%
20	15.839	2180	2185	2191	rVB5	38683	87118	0.37%	0.037%
21	16.051	2216	2221	2227	rBV2	193162	363258	1.55%	0.156%
22	16.663	2320	2325	2329	rVB2	66154	114219	0.49%	0.049%
23	16.739	2334	2338	2345	rVB4	65685	109897	0.47%	0.047%
24	16.816	2345	2351	2355	rBV4	54150	115449	0.49%	0.049%
25	16.863	2355	2359	2363	rVV3	82895	120969	0.52%	0.052%
26	16.910	2363	2367	2374	rVB4	52341	108758	0.46%	0.047%
27	17.081	2389	2396	2401	rBV2	5561029	8178321	34.94%	3.502%
28	17.122	2401	2403	2408	rVB	591336	726842	3.11%	0.311%
29	17.181	2408	2413	2415	rBV	317206	458862	1.96%	0.197%
30	17.216	2415	2419	2424	rVV	1081139	1505923	6.43%	0.645%
31	17.269	2424	2428	2436	rVB	146000	215417	0.92%	0.092%
32	17.486	2457	2465	2468	rBV	261324	407693	1.74%	0.175%
33	17.528	2468	2472	2482	rVV	141318	349237	1.49%	0.150%
34	17.604	2482	2485	2492	rVB2	180139	250572	1.07%	0.107%

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Integration Parameters: LSCINT.P

Integrator: RTE
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 Title : SVOA CALIBRATION

35	17.957	2538	2545	2548	rBV2	139518	266504	1.14%	0.114%
36	18.004	2548	2553	2561	rVB	280330	516692	2.21%	0.221%
37	18.086	2561	2567	2571	rBV	388756	600472	2.57%	0.257%
38	18.151	2572	2578	2582	rBV3	311963	677864	2.90%	0.290%
39	18.275	2596	2599	2604	rBV3	68992	127590	0.55%	0.055%
40	18.351	2610	2612	2617	rVB4	84885	96100	0.41%	0.041%
41	18.463	2627	2631	2634	rBV	177876	225163	0.96%	0.096%
42	18.498	2634	2637	2643	rVB3	200173	281455	1.20%	0.121%
43	18.586	2650	2652	2656	rBV3	65664	83763	0.36%	0.036%
44	18.710	2665	2673	2677	rBV3	243381	493055	2.11%	0.211%
45	18.763	2679	2682	2685	rVV2	204064	282207	1.21%	0.121%
46	18.798	2685	2688	2693	rVB	261751	354494	1.51%	0.152%
47	18.880	2697	2702	2708	rBV	485975	801188	3.42%	0.343%
48	18.939	2708	2712	2716	rVV2	272650	462240	1.98%	0.198%
49	18.975	2716	2718	2721	rVB	167644	159542	0.68%	0.068%
50	19.022	2721	2726	2735	rVB	1017450	1429797	6.11%	0.612%
51	19.145	2741	2747	2754	rVB	10578418	14044031	60.01%	6.014%
52	19.216	2755	2759	2761	rBV	251515	331855	1.42%	0.142%
53	19.245	2761	2764	2768	rVB3	248224	324970	1.39%	0.139%
54	19.292	2769	2772	2776	rVB	238232	309559	1.32%	0.133%
55	19.339	2776	2780	2783	rBV5	126983	224483	0.96%	0.096%
56	19.392	2786	2789	2792	rVB	231213	281453	1.20%	0.121%
57	19.510	2799	2809	2817	rBV	13474013	22031631	94.13%	9.435%
58	19.586	2817	2822	2830	rVV4	501103	1069438	4.57%	0.458%
59	19.692	2836	2840	2845	rBV	488605	676177	2.89%	0.290%
60	19.751	2846	2850	2856	rVB2	799890	1266733	5.41%	0.542%
61	19.839	2859	2865	2876	rBV4	1492671	3811093	16.28%	1.632%
62	19.927	2877	2880	2883	rBV3	179954	207311	0.89%	0.089%
63	19.980	2883	2889	2898	rVB3	1556281	4919806	21.02%	2.107%
64	20.057	2898	2902	2905	rBV2	242609	371801	1.59%	0.159%
65	20.110	2907	2911	2914	rVB	625342	748787	3.20%	0.321%
66	20.169	2914	2921	2925	rBV2	2009579	3571814	15.26%	1.530%
67	20.210	2925	2928	2931	rVB	436523	498874	2.13%	0.214%
68	20.245	2931	2934	2940	rVB2	323063	492729	2.11%	0.211%
69	20.310	2940	2945	2948	rBV	1464977	2042437	8.73%	0.875%
70	20.357	2948	2953	2957	rVB2	967270	1453462	6.21%	0.622%
71	20.569	2985	2989	2991	rBV3	440540	709276	3.03%	0.304%

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Integration Parameters: LSCINT.P

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72	20.674	3005	3007	3010	rVB3	381398	352237	1.50%	0.151%
73	20.757	3017	3021	3028	rVB	2120042	2967959	12.68%	1.271%
74	20.927	3043	3050	3053	rBV3	1735069	3217348	13.75%	1.378%
75	20.969	3053	3057	3059	rVV	1487006	1884453	8.05%	0.807%
76	21.004	3059	3063	3068	rVV2	1972133	3048079	13.02%	1.305%
77	21.057	3068	3072	3080	rVB2	1417382	2133238	9.11%	0.914%
78	21.169	3084	3091	3096	rBV4	629362	1474132	6.30%	0.631%
79	21.269	3099	3108	3114	rBV2	9987628	23404524	100.00%	10.023%
80	21.322	3114	3117	3124	rVV	7966138	10621219	45.38%	4.548%
81	21.398	3128	3130	3133	rVB3	294387	288156	1.23%	0.123%
82	21.439	3133	3137	3142	rBV2	582928	1163067	4.97%	0.498%
83	21.592	3159	3163	3168	rBV2	1182011	1835276	7.84%	0.786%
84	21.645	3168	3172	3175	rBV3	692560	933560	3.99%	0.400%
85	21.780	3192	3195	3198	rBV	459111	592800	2.53%	0.254%
86	21.833	3198	3204	3208	rVV	1787159	3063618	13.09%	1.312%
87	21.886	3209	3213	3219	rVB2	1295681	1921288	8.21%	0.823%
88	22.027	3233	3237	3241	rBV	862854	1296791	5.54%	0.555%
89	22.133	3251	3255	3260	rVB2	580743	789649	3.37%	0.338%
90	22.304	3281	3284	3288	rBV	409297	518387	2.21%	0.222%
91	22.598	3327	3334	3345	rVB3	859090	2569478	10.98%	1.100%
92	22.857	3370	3378	3382	rBV	9728077	21490734	91.82%	9.203%
93	23.021	3401	3406	3412	rVB	1328948	2132792	9.11%	0.913%
94	23.151	3423	3428	3437	rBV	695941	1880483	8.03%	0.805%
95	23.327	3451	3458	3468	rVB	5126282	9845495	42.07%	4.216%
96	23.421	3468	3474	3481	rVB	5531479	9874706	42.19%	4.229%
97	23.515	3483	3490	3494	rVV	2751042	5657248	24.17%	2.423%
98	23.563	3494	3498	3506	rVB2	1891115	3994081	17.07%	1.710%
99	25.751	3861	3870	3882	rVB2	3104084	9053802	38.68%	3.877%
100	26.427	3976	3985	4002	rVB	1932803	6335080	27.07%	2.713%

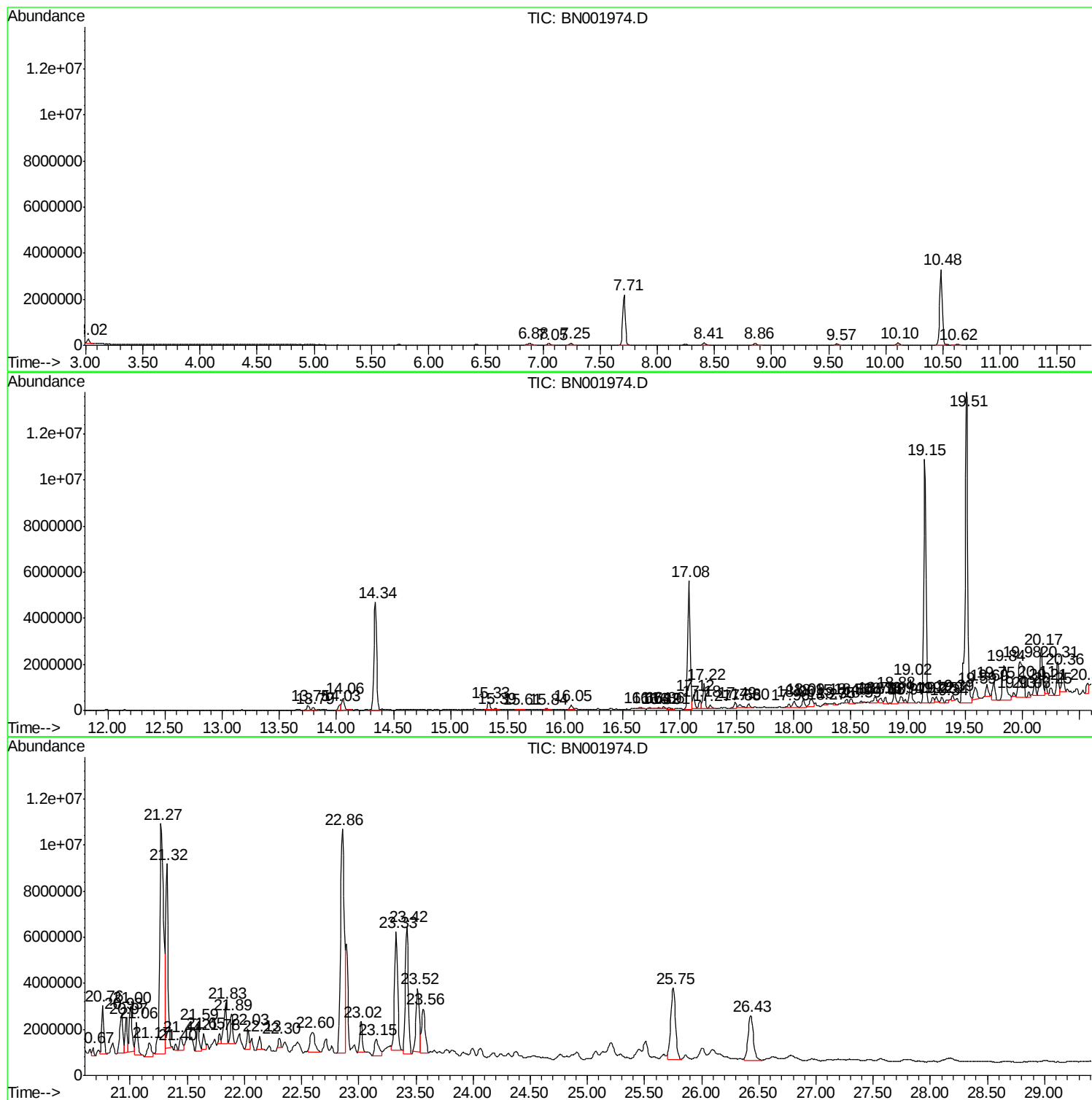
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Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061918.M
Quant Title : SVOA CALIBRATION

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



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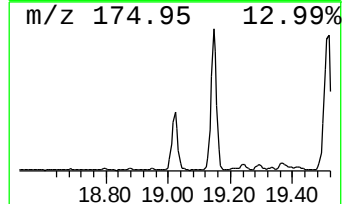
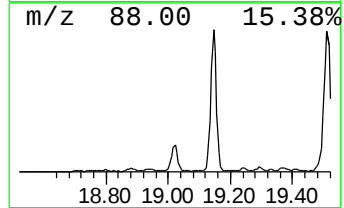
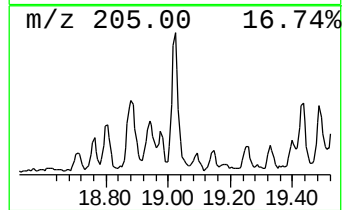
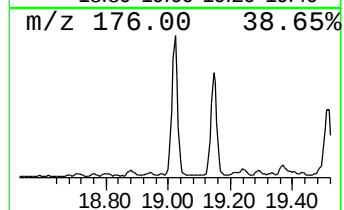
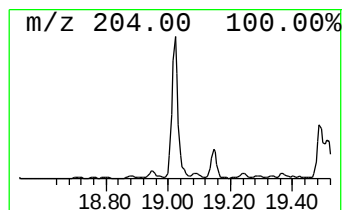
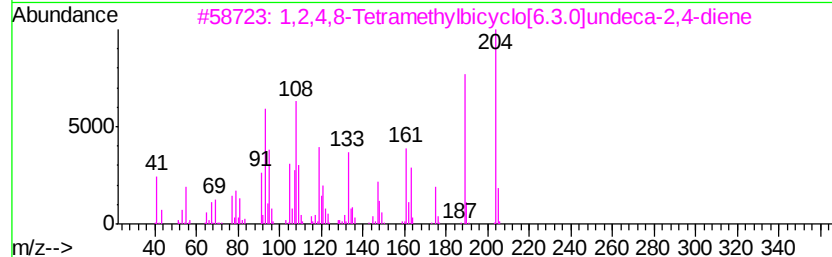
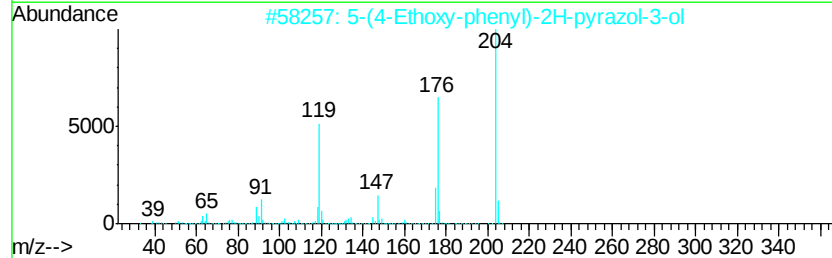
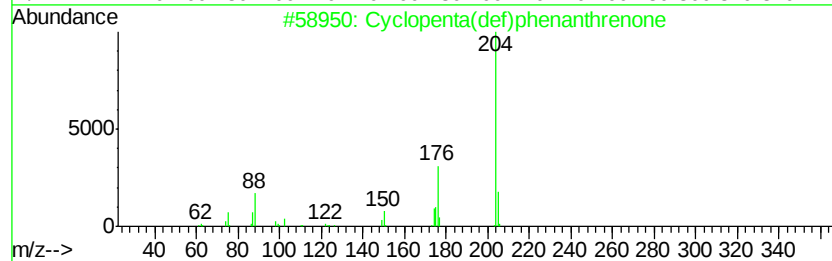
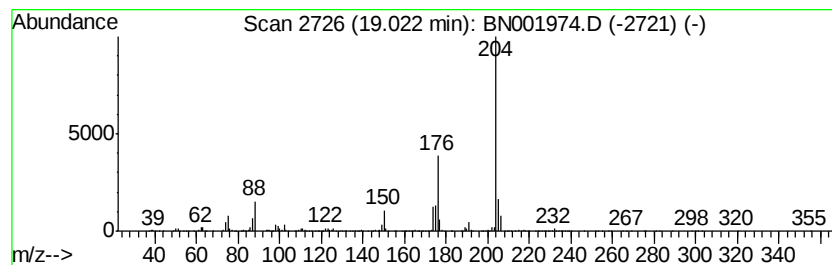
Instrument :
BNA_N
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F1H00DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061918.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Cyclopenta(def)phenanthrenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.02	3.50 ng/ul	1429800	Phenanthrene-d10	17.08	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	97
2	5-(4-Ethoxy-phenyl)-2H-pyrazol-3-ol	204	C11H12N2O2	1000278-26-8	59
3	1,2,4,8-Tetramethylbicyclo[6.3.0...	204	C15H24	137235-51-9	43
4	Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	43
5	4(equat)-(but-3-en-1-ynyl)-2(axi...	219	C14H21NO	038423-22-2	42



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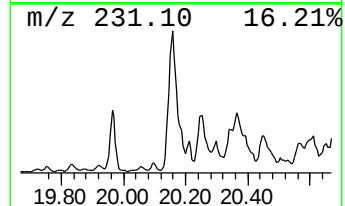
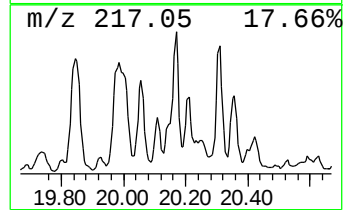
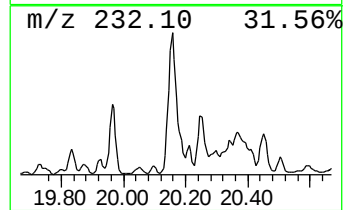
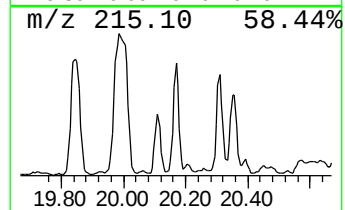
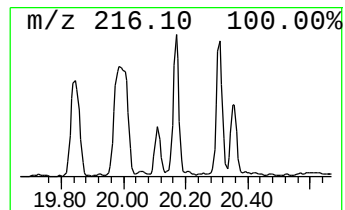
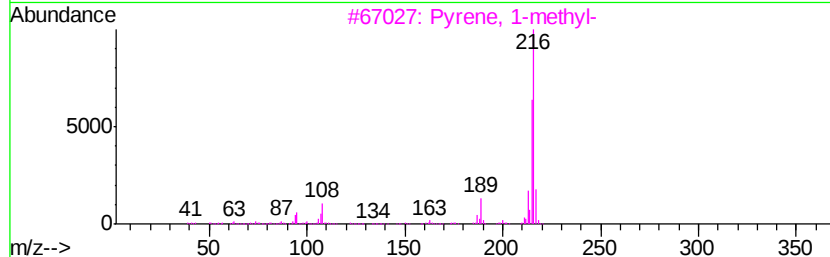
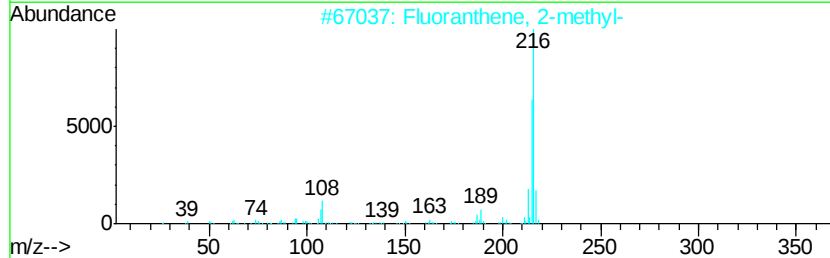
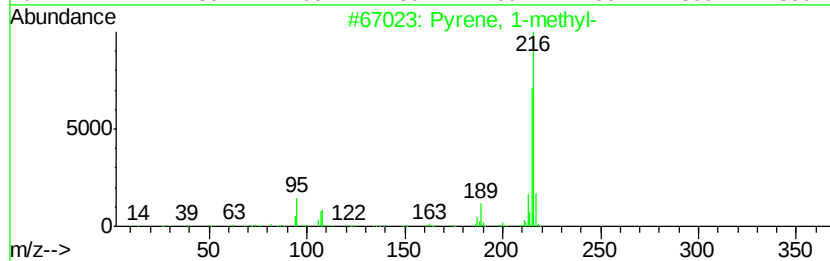
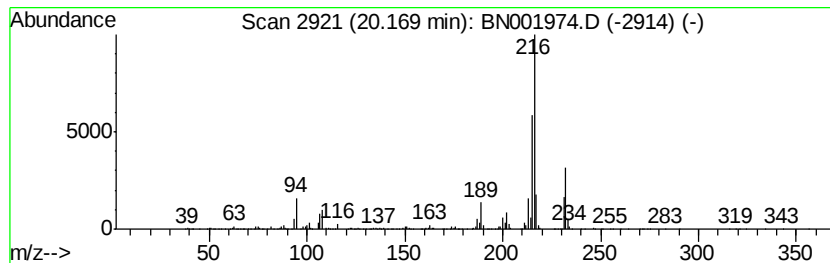
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061918.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Pyrene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.17	3.05 ng/ul	3571810	Chrysene-d12	21.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyrene, 1-methyl-	216	C17H12	002381-21-7	98
2		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	98
3		Pyrene, 1-methyl-	216	C17H12	002381-21-7	96
4		Pyrene, 4-methyl-	216	C17H12	003353-12-6	96
5		Pyrene, 2-methyl-	216	C17H12	003442-78-2	95



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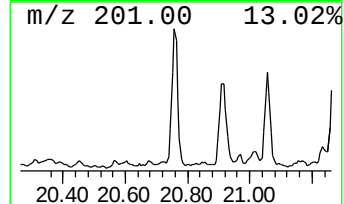
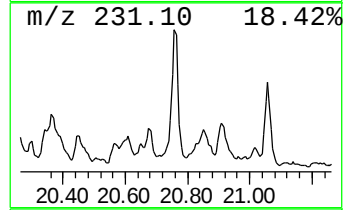
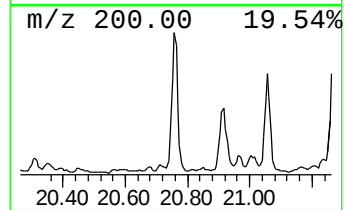
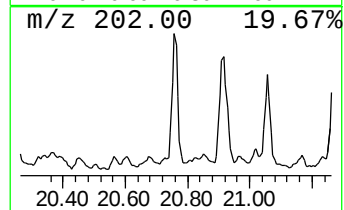
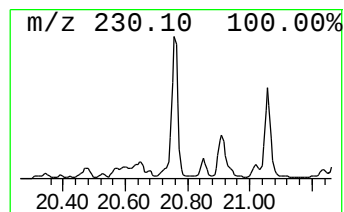
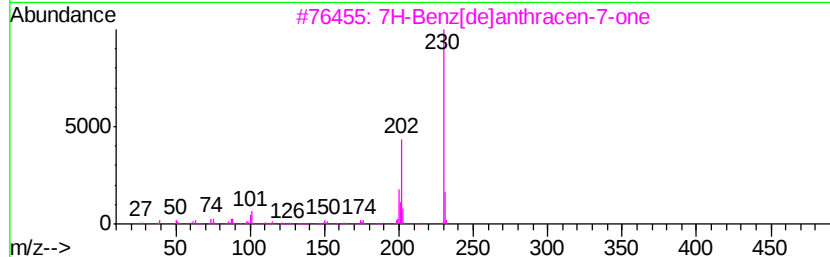
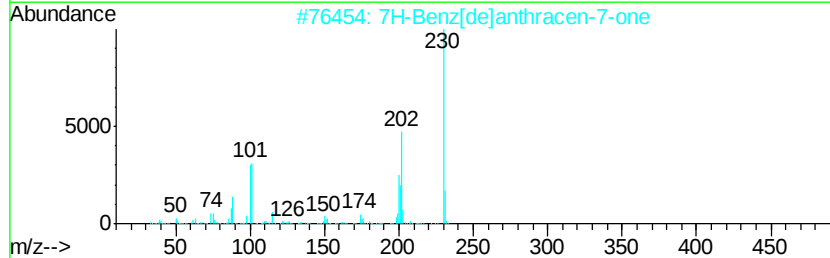
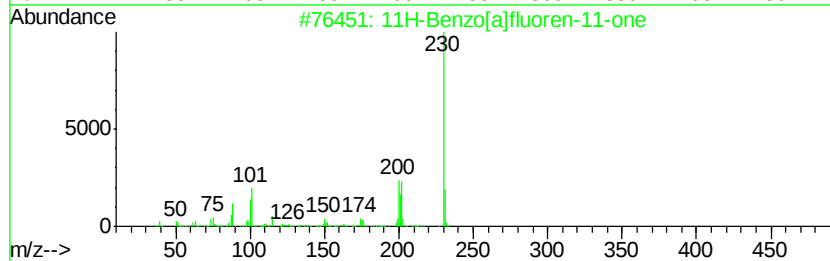
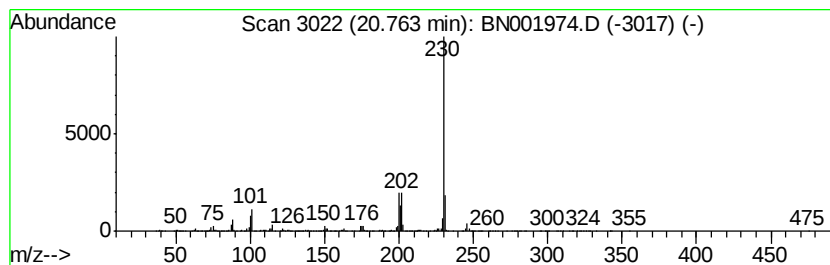
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.76	2.54 ng/ul	2967960	Chrysene-d12	21.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	95
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	90
4		7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	62
5		7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	59



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN070518\
Data File : BN001974.D
Acq On : 05 Jul 2018 19:06
Operator : JU/SJ
Sample : J3721-14DL 25X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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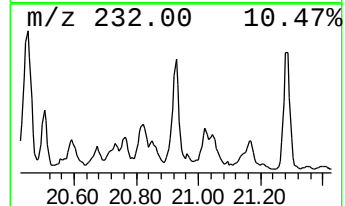
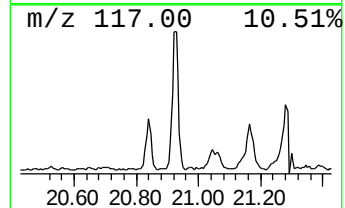
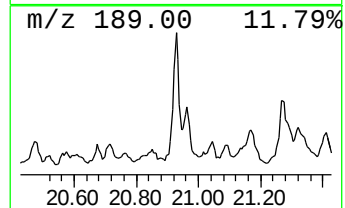
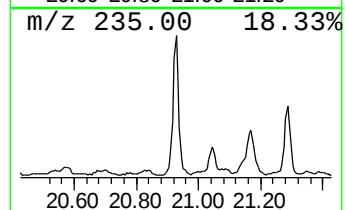
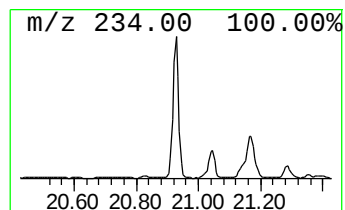
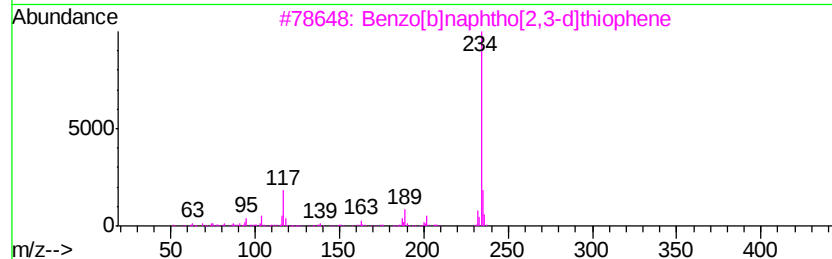
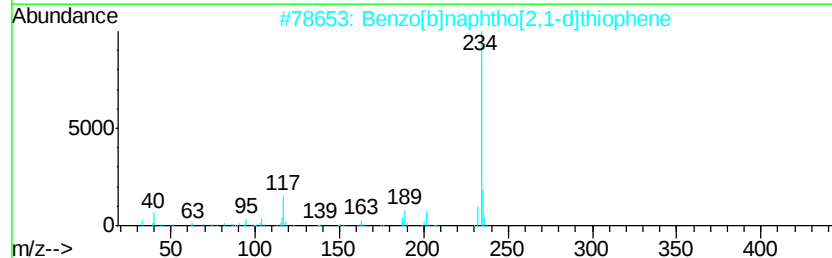
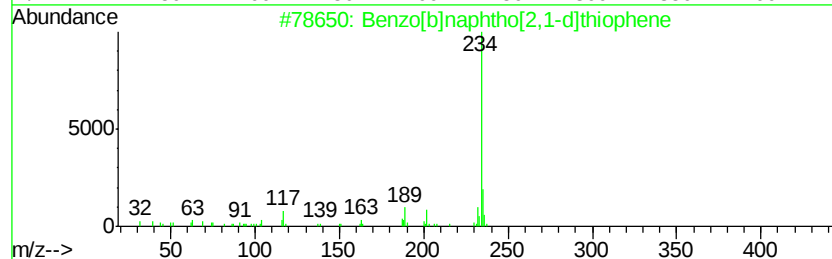
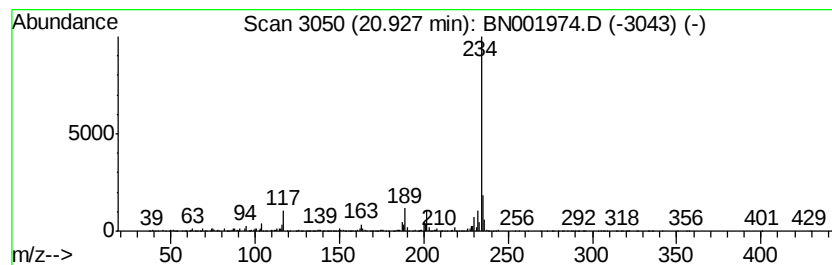
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzo[b]naphtho[2,1-d]thiop... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.93	2.75 ng/ul	3217350	Chrysene-d12	21.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	96
2		Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	94
3		Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	94
4		Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	91
5		Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	91



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN070518\
Data File : BN001974.D
Acq On : 05 Jul 2018 19:06
Operator : JU/SJ
Sample : J3721-14DL 25X
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Instrument :
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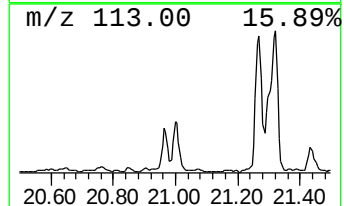
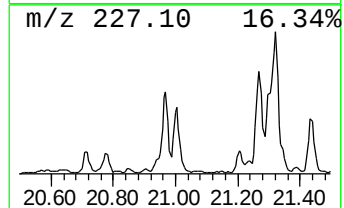
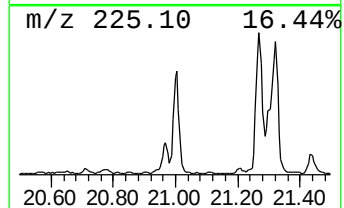
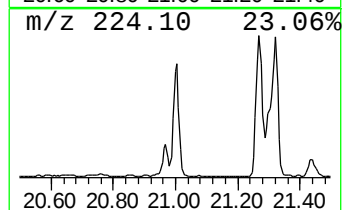
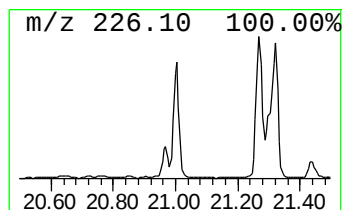
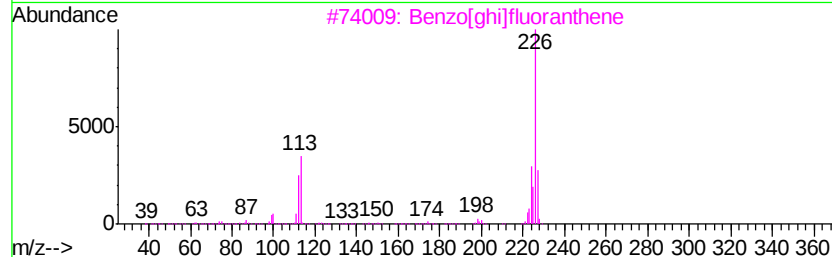
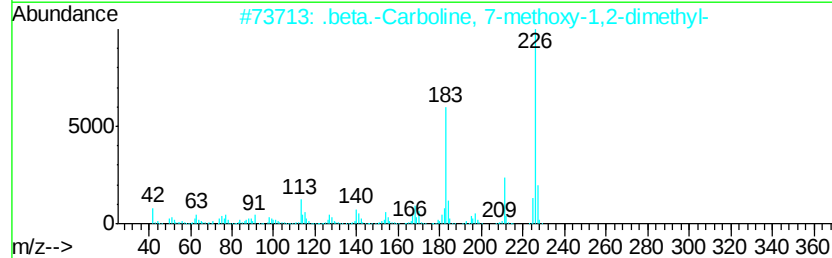
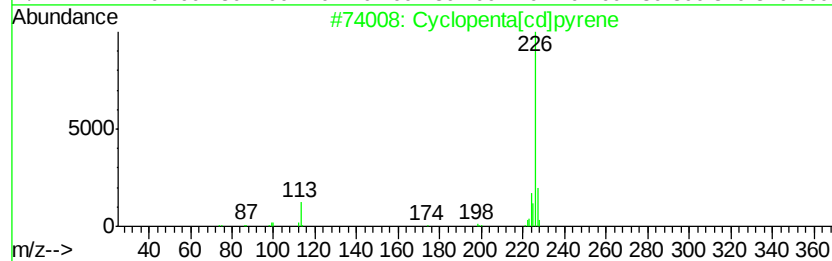
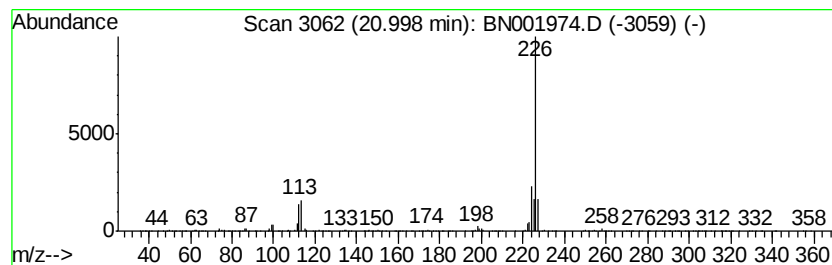
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Cyclopenta[cd]pyrene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.00	2.60 ng/ul	3048080	Chrysene-d12	21.29

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	76
2		.beta.-Carboline, 7-methoxy-1,2-...	226	C14H14N2O	006519-18-2	53
3		Benzo[ghi]fluoranthene	226	C18H10	000203-12-3	43
4		9H-Xanthen-9-one, 3-methoxy-	226	C14H10O3	003722-52-9	38
5		1,3-Diamino-5,6-dihydro-9-methyl...	226	C13H14N4	037436-39-8	36



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN070518\
Data File : BN001974.D
Acq On : 05 Jul 2018 19:06
Operator : JU/SJ
Sample : J3721-14DL 25X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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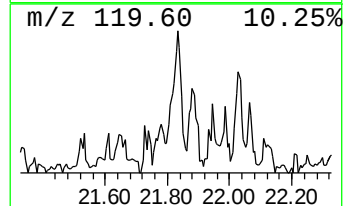
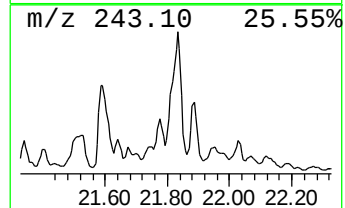
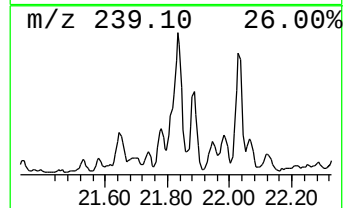
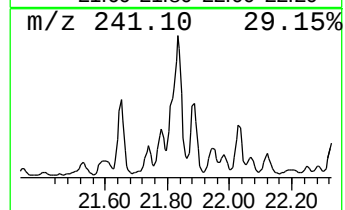
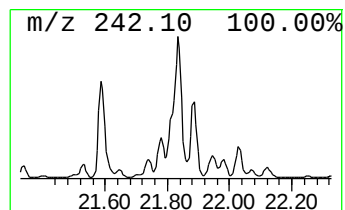
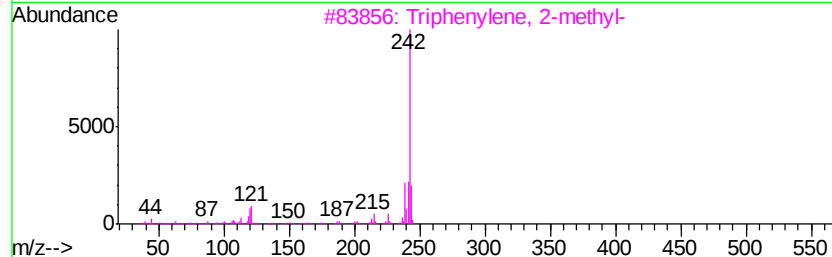
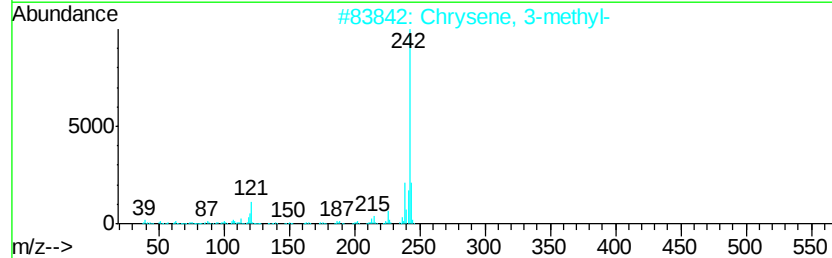
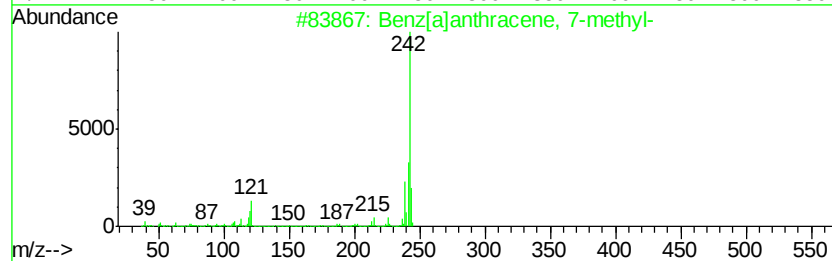
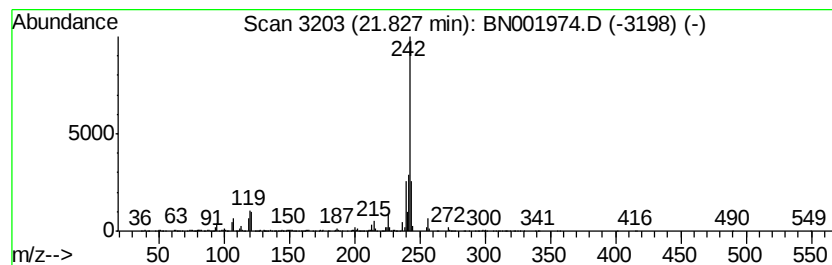
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benz[a]anthracene, 7-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.83	2.62 ng/ul	3063620	Chrysene-d12	21.29

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benz[a]anthracene, 7-methyl-	242	C19H14	002541-69-7	94
2			Chrysene, 3-methyl-	242	C19H14	003351-31-3	93
3			Triphenylene, 2-methyl-	242	C19H14	001705-84-6	93
4			Benz[a]anthracene, 7-methyl-	242	C19H14	002541-69-7	93
5			Benz[a]anthracene, 1-methyl-	242	C19H14	002498-77-3	89



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN070518\
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Acq On : 05 Jul 2018 19:06
Operator : JU/SJ
Sample : J3721-14DL 25X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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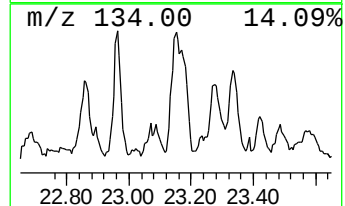
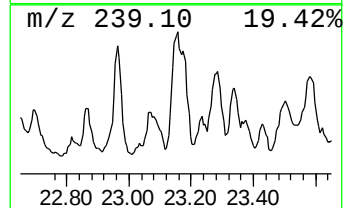
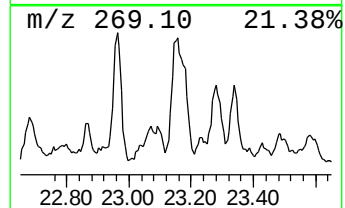
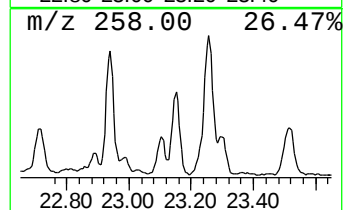
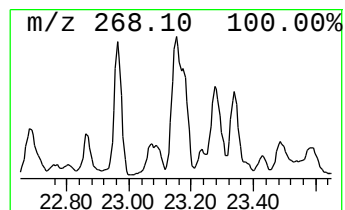
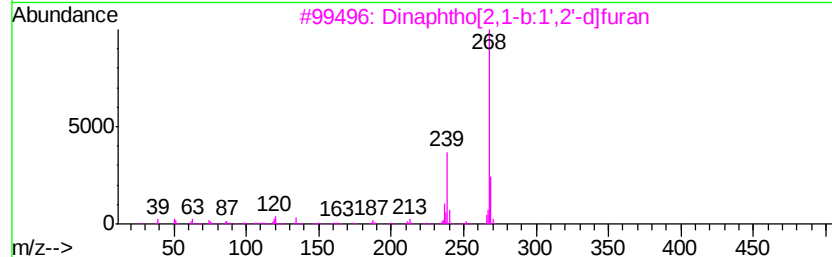
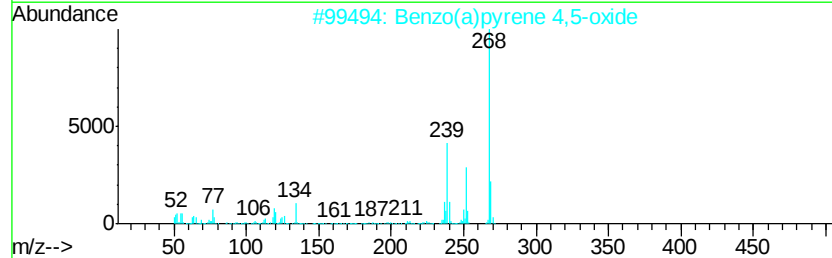
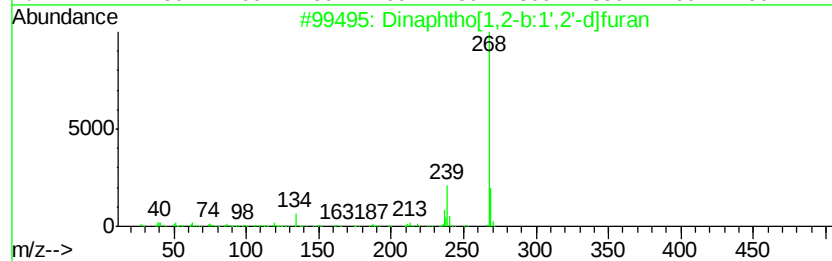
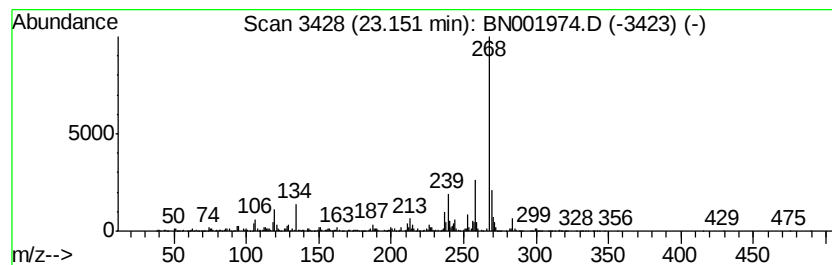
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Dinaphtho[1,2-b:1',2'-d]furan Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.15	6.65 ng/ul	1880480	Perylene-d12	23.52

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dinaphtho[1,2-b:1',2'-d]furan	268	C20H12O	000207-93-2	81
2		Benzo(a)pyrene 4,5-oxide	268	C20H12O	037574-47-3	72
3		Dinaphtho[2,1-b:1',2'-d]furan	268	C20H12O	000194-63-8	68
4		10-Hydroxy-5-methoxy-2-methyl-1,...	268	C16H12O4	084542-45-0	58
5		9,10-Anthracenedione, 1,5-dimeth...	268	C16H12O4	006448-90-4	53



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN070518\
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Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN061918.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
Cyclopenta(def)ph...	19.02	3.5	ng/ul	1429800	4	17.08	8178320	20.0
Pyrene, 1-methyl-	20.17	3.0	ng/ul	3571810	5	21.29	23404500	20.0
11H-Benzo[a]fluor...	20.76	2.5	ng/ul	2967960	5	21.29	23404500	20.0
Benzo[b]naphtho[2...	20.93	2.8	ng/ul	3217350	5	21.29	23404500	20.0
Cyclopenta[cd]pyrene	21.00	2.6	ng/ul	3048080	5	21.29	23404500	20.0
Benz[a]anthracene...	21.83	2.6	ng/ul	3063620	5	21.29	23404500	20.0
Dinaphtho[1,2-b:1...	23.15	6.7	ng/ul	1880480	6	23.52	5657250	20.0