

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030320\
 Data File : BN010075.D
 Acq On : 03 Mar 2020 23:22
 Operator : CG/JU
 Sample : L1619-04DL 10X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBDT8DL

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-BN021220MA.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	3.511	86	89	93	rVB6	5094	7625	0.28%	0.039%
2	3.687	118	119	123	rBV4	3084	3444	0.13%	0.018%
3	4.234	207	212	217	rBV5	4546	8201	0.30%	0.042%
4	6.617	614	617	620	rBV2	3730	5353	0.20%	0.027%
5	6.758	635	641	646	rBV3	4934	9296	0.34%	0.047%
6	6.928	667	670	676	rBV4	6153	10053	0.37%	0.051%
7	7.369	738	745	762	rBV	209679	378144	13.86%	1.932%
8	7.916	832	838	846	rBV3	9193	19277	0.71%	0.098%
9	8.546	943	945	952	rBV4	4376	8451	0.31%	0.043%
10	10.116	1203	1212	1228	rBV	212421	424752	15.57%	2.170%
11	11.828	1496	1503	1510	rBV3	2991	6484	0.24%	0.033%
12	12.928	1687	1690	1695	rVB2	2819	3551	0.13%	0.018%
13	13.457	1775	1780	1785	rBV2	19143	32232	1.18%	0.165%
14	13.499	1785	1787	1794	rVB3	3568	5307	0.19%	0.027%
15	13.698	1816	1821	1824	rBV	20804	34049	1.25%	0.174%
16	13.728	1824	1826	1836	rVB3	19421	32410	1.19%	0.166%
17	14.004	1868	1873	1883	rBV	437225	673293	24.68%	3.440%
18	14.069	1883	1884	1890	rVB4	3502	4794	0.18%	0.024%
19	14.440	1942	1947	1952	rVB	9889	14991	0.55%	0.077%
20	15.022	2041	2046	2051	rBV	31945	45954	1.68%	0.235%
21	15.081	2051	2056	2065	rVB2	31794	52968	1.94%	0.271%
22	15.316	2092	2096	2097	rBV2	3289	3986	0.15%	0.020%
23	15.787	2170	2176	2180	rBV3	12313	26059	0.96%	0.133%
24	16.098	2222	2229	2231	rBV3	3163	4561	0.17%	0.023%
25	16.563	2304	2308	2310	rBV2	6909	8004	0.29%	0.041%
26	16.616	2312	2317	2322	rVB5	11834	19040	0.70%	0.097%
27	16.769	2337	2343	2347	rBV	616014	861465	31.58%	4.401%
28	16.810	2347	2350	2355	rVB	88328	114053	4.18%	0.583%
29	16.898	2360	2365	2376	rVV	1044044	1513336	55.48%	7.731%
30	16.981	2376	2379	2385	rVV3	6062	12287	0.45%	0.063%
31	17.051	2389	2391	2395	rVB2	4412	4875	0.18%	0.025%
32	17.187	2409	2414	2425	rBV	165096	264698	9.70%	1.352%
33	17.298	2428	2433	2436	rVV4	6728	10558	0.39%	0.054%
34	17.698	2497	2501	2504	rBV4	7956	11033	0.40%	0.056%

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030320\
 Data File : BN010075.D
 Acq On : 03 Mar 2020 23:22
 Operator : CG/JU
 Sample : L1619-04DL 10X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBDT8DL

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-BN021220MA.M
 Title : SVOA CALIBRATION

35	17.787	2512	2516	2522	rBV	24349	34405	1.26%	0.176%
36	17.851	2522	2527	2530	rBV5	9789	16588	0.61%	0.085%
37	17.881	2530	2532	2535	rVB4	4786	4787	0.18%	0.024%
38	17.975	2543	2548	2549	rBV2	6127	9130	0.33%	0.047%
39	18.010	2553	2554	2558	rVB2	8442	6220	0.23%	0.032%
40	18.051	2558	2561	2565	rBV2	21840	35625	1.31%	0.182%
41	18.128	2571	2574	2578	rVB3	13272	18891	0.69%	0.097%
42	18.163	2578	2580	2583	rVB3	4170	3832	0.14%	0.020%
43	18.216	2583	2589	2596	rBV6	32718	64798	2.38%	0.331%
44	18.457	2628	2630	2632	rVB2	8182	6088	0.22%	0.031%
45	18.492	2632	2636	2640	rBV5	10548	16429	0.60%	0.084%
46	18.575	2646	2650	2654	rBV6	14083	22818	0.84%	0.117%
47	18.634	2657	2660	2662	rBV3	10121	11215	0.41%	0.057%
48	18.681	2667	2668	2670	rVB2	10489	5566	0.20%	0.028%
49	18.734	2670	2677	2679	rBV2	26418	41965	1.54%	0.214%
50	18.839	2690	2695	2701	rVV	157241	215383	7.90%	1.100%
51	18.916	2706	2708	2709	rVV2	9596	7260	0.27%	0.037%
52	18.951	2710	2714	2718	rVV2	25322	44046	1.61%	0.225%
53	19.004	2720	2723	2726	rVV4	20494	25797	0.95%	0.132%
54	19.045	2729	2730	2731	rVV	6028	4009	0.15%	0.020%
55	19.081	2732	2736	2740	rVV2	40124	58231	2.13%	0.297%
56	19.128	2742	2744	2748	rVB5	15567	16610	0.61%	0.085%
57	19.181	2748	2753	2754	rBV3	85656	117163	4.30%	0.599%
58	19.204	2754	2757	2762	rVV	206006	289013	10.59%	1.476%
59	19.239	2762	2763	2769	rVV5	25625	32274	1.18%	0.165%
60	19.292	2769	2772	2779	rVB8	20463	34463	1.26%	0.176%
61	19.398	2787	2790	2795	rBV4	22019	29919	1.10%	0.153%
62	19.463	2797	2801	2806	rVB	50947	77621	2.85%	0.397%
63	19.551	2810	2816	2823	rBV4	39434	98409	3.61%	0.503%
64	19.692	2833	2840	2848	rBV3	101526	226501	8.30%	1.157%
65	19.769	2850	2853	2854	rBV2	11032	10127	0.37%	0.052%
66	19.816	2858	2861	2865	rVB5	16116	18846	0.69%	0.096%
67	19.875	2865	2871	2875	rBV3	69359	136224	4.99%	0.696%
68	20.016	2890	2895	2899	rBV	81656	126927	4.65%	0.648%
69	20.057	2899	2902	2908	rVV4	32300	60439	2.22%	0.309%
70	20.269	2935	2938	2942	rBV6	24739	38248	1.40%	0.195%
71	20.375	2954	2956	2961	rBV5	24514	34259	1.26%	0.175%

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN030320\
 Data File : BN010075.D
 Acq On : 03 Mar 2020 23:22
 Operator : CG/JU
 Sample : L1619-04DL 10X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DBDT8DL

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-BN021220MA.M
 Title : SVOA CALIBRATION

72	20.433	2961	2966	2969	rBV3	44526	81955	3.00%	0.419%
73	20.475	2969	2973	2978	rBV5	55365	103243	3.78%	0.527%
74	20.633	2997	3000	3003	rBV	43001	55286	2.03%	0.282%
75	20.669	3003	3006	3009	rBV3	70658	96840	3.55%	0.495%
76	20.704	3009	3012	3016	rBV	189162	224575	8.23%	1.147%
77	20.780	3022	3025	3029	rBV2	89706	126632	4.64%	0.647%
78	20.810	3029	3030	3034	rVB	60896	58704	2.15%	0.300%
79	20.986	3050	3060	3064	rBV3	1026618	2067638	75.80%	10.563%
80	21.022	3064	3066	3072	rVB	342502	399782	14.66%	2.042%
81	21.110	3078	3081	3084	rBV5	41900	54181	1.99%	0.277%
82	21.292	3109	3112	3119	rVB	118661	174061	6.38%	0.889%
83	21.357	3119	3123	3126	rBV3	65344	100615	3.69%	0.514%
84	21.475	3140	3143	3146	rBV	46757	66508	2.44%	0.340%
85	21.528	3149	3152	3155	rVV	140547	178992	6.56%	0.914%
86	21.575	3158	3160	3166	rVB3	70352	117621	4.31%	0.601%
87	21.804	3197	3199	3203	rVB	56201	65360	2.40%	0.334%
88	22.239	3269	3273	3284	rVB4	160570	294351	10.79%	1.504%
89	22.357	3290	3293	3297	rVB2	81073	101042	3.70%	0.516%
90	22.475	3306	3313	3317	rBV2	1339396	2727880	100.00%	13.936%
91	22.627	3335	3339	3344	rVB	133730	202001	7.41%	1.032%
92	22.904	3380	3386	3391	rBV	581195	954160	34.98%	4.874%
93	22.992	3396	3401	3408	rVB	569714	924002	33.87%	4.720%
94	23.074	3410	3415	3420	rVV2	703854	1210217	44.36%	6.183%
95	23.122	3420	3423	3434	rVB3	176403	381738	13.99%	1.950%
96	23.569	3495	3499	3506	rVB2	96364	159590	5.85%	0.815%
97	24.233	3608	3612	3618	rVB2	81042	141335	5.18%	0.722%
98	24.698	3688	3691	3703	rVB3	110739	271570	9.96%	1.387%
99	25.110	3754	3761	3771	rVB2	299850	773263	28.35%	3.950%
100	25.339	3794	3800	3810	rVB2	375795	826810	30.31%	4.224%

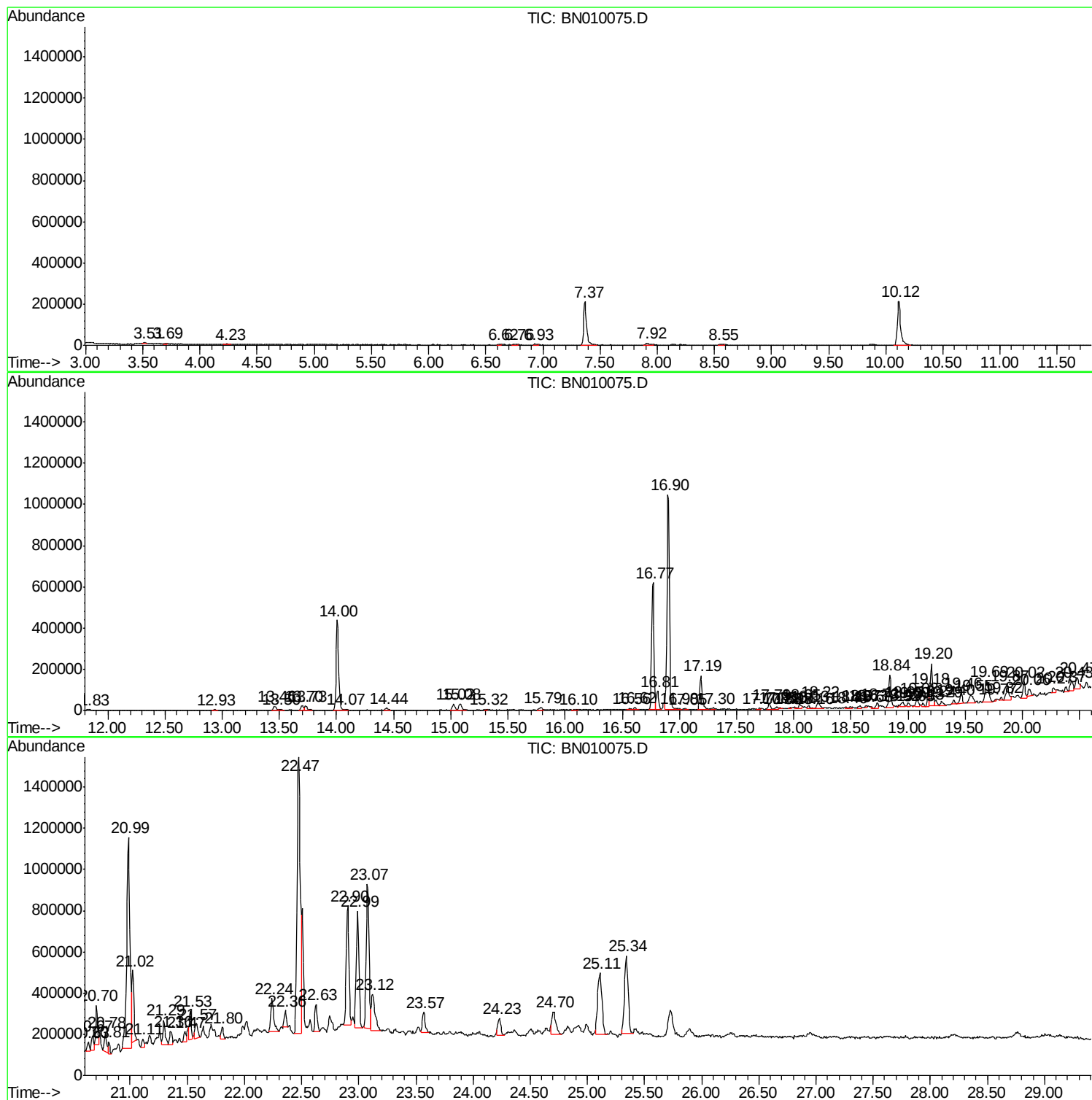
Sum of corrected areas: 19574662

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030320\
Data File : BN010075.D
Acq On : 03 Mar 2020 23:22
Operator : CG/JU
Sample : L1619-04DL 10X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBDT8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030320\
Data File : BN010075.D
Acq On : 03 Mar 2020 23:22
Operator : CG/JU
Sample : L1619-04DL 10X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBDT8DL

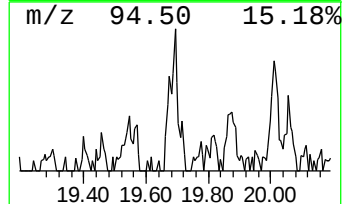
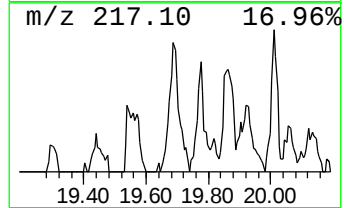
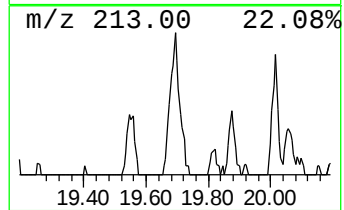
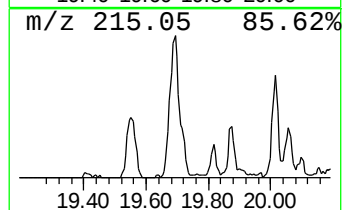
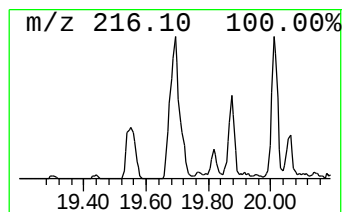
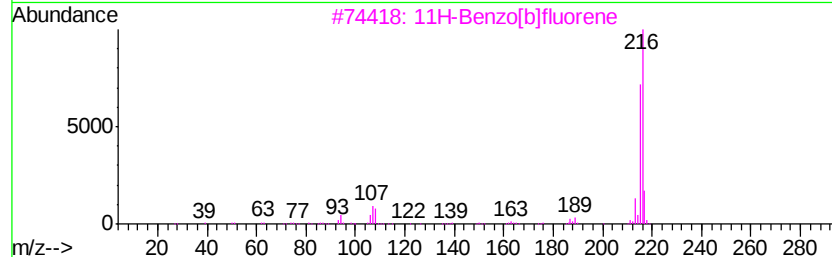
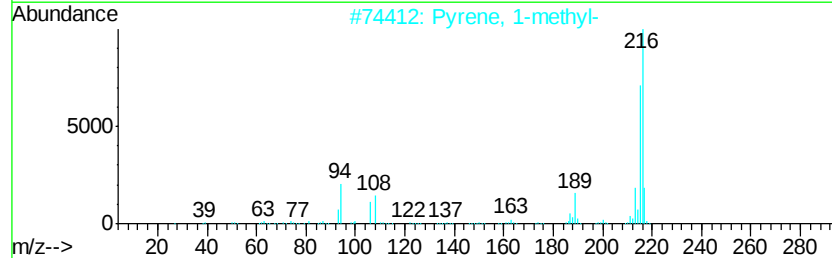
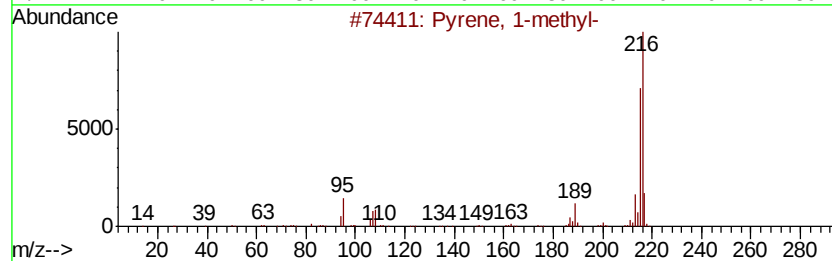
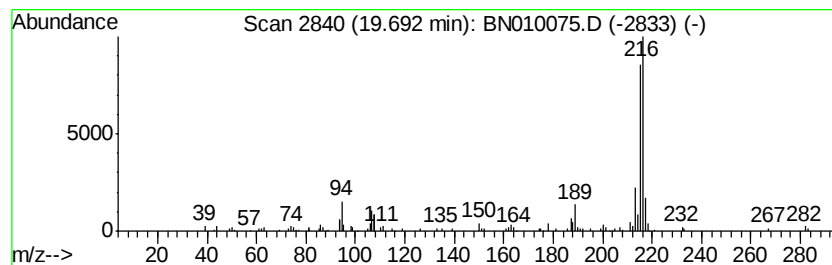
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.69	2.19 ng/ul	226501	Chrysene-d12	20.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
2		Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
3		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
4		Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
5		11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030320\
Data File : BN010075.D
Acq On : 03 Mar 2020 23:22
Operator : CG/JU
Sample : L1619-04DL 10X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBDT8DL

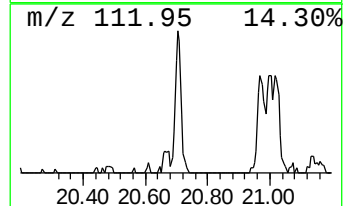
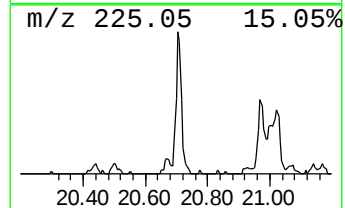
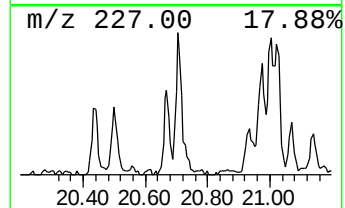
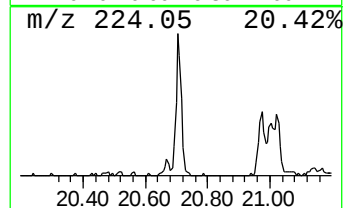
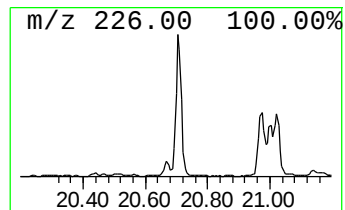
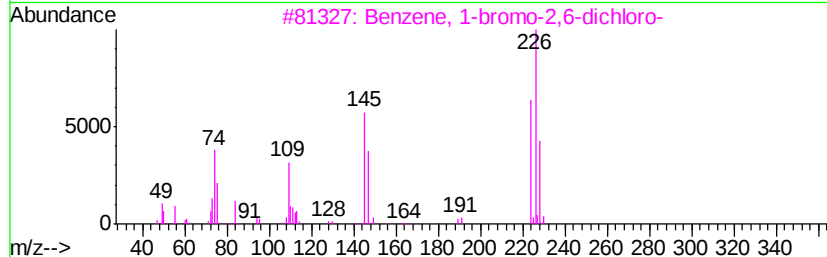
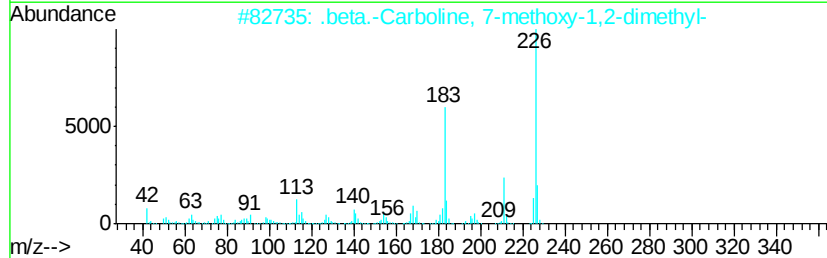
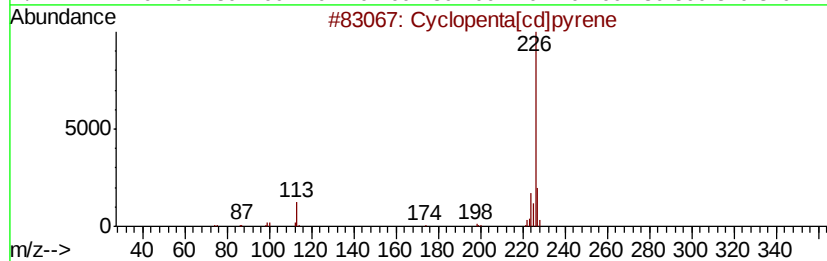
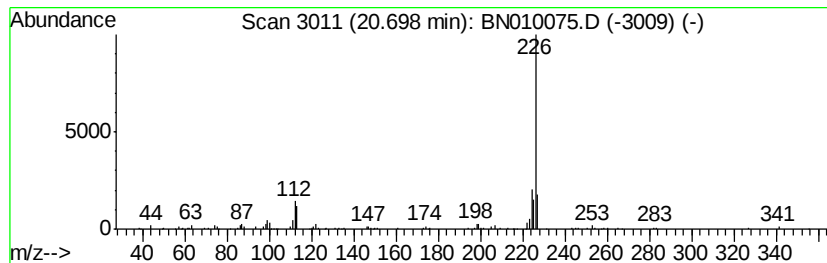
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Cyclopenta[cd]pyrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.70	2.17 ng/ul	224575	Chrysene-d12	20.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta[cd]pyrene	226	C18H10	027208-37-3	74
2		.beta.-Carboline, 7-methoxy-1,2-...	226	C14H14N2O	006519-18-2	53
3		Benzene, 1-bromo-2,6-dichloro-	224	C6H3BrCl2	019393-92-1	50
4		Benzo[ghi]fluoranthene	226	C18H10	000203-12-3	47
5		1,8-Diazacyclotetradecane-2,7-dione	226	C12H22N2O2	004266-66-4	37



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030320\
Data File : BN010075.D
Acq On : 03 Mar 2020 23:22
Operator : CG/JU
Sample : L1619-04DL 10X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBDT8DL

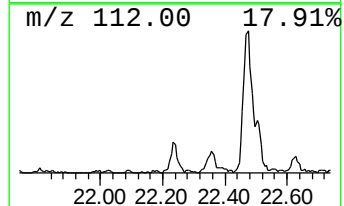
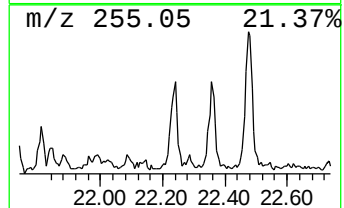
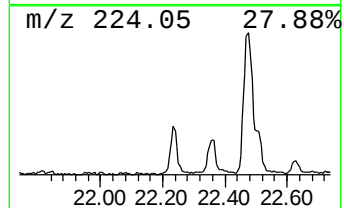
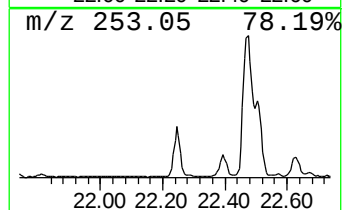
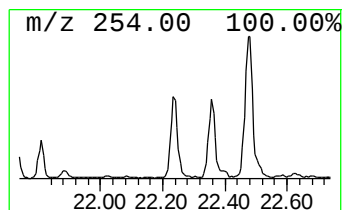
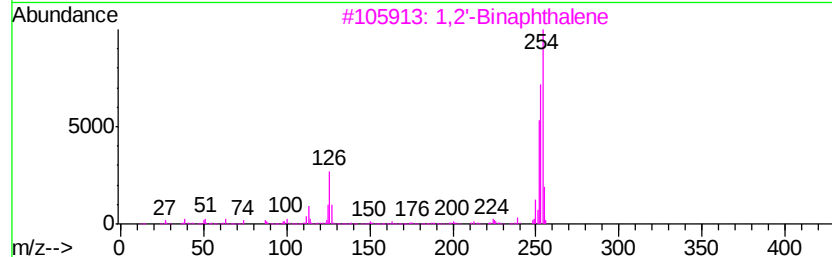
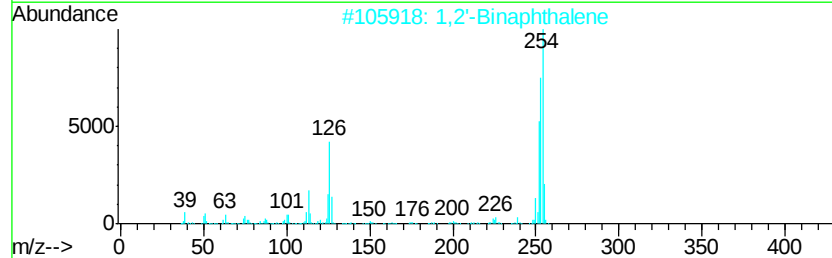
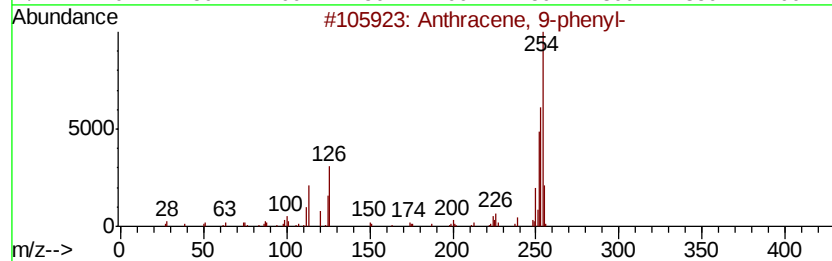
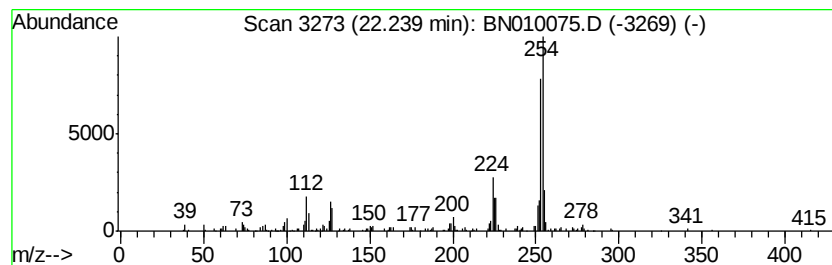
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Anthracene, 9-phenyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.24	4.86 ng/ul	294351	Perylene-d12	23.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Anthracene, 9-phenyl-	254	C20H14	000602-55-1	74
2		1,2'-Binaphthalene	254	C20H14	004325-74-0	72
3		1,2'-Binaphthalene	254	C20H14	004325-74-0	68
4		1,1'-Binaphthalene	254	C20H14	000604-53-5	68
5		1,1'-Binaphthalene	254	C20H14	000604-53-5	64



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030320\
Data File : BN010075.D
Acq On : 03 Mar 2020 23:22
Operator : CG/JU
Sample : L1619-04DL 10X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBDT8DL

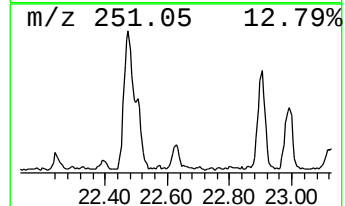
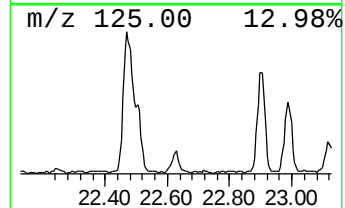
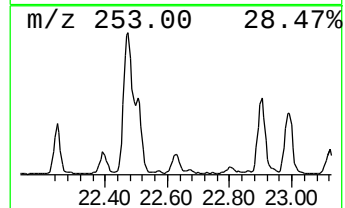
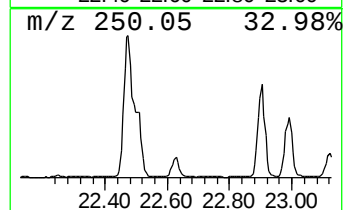
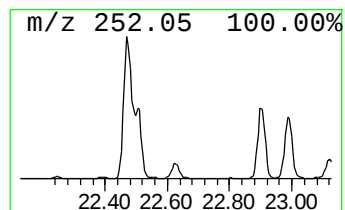
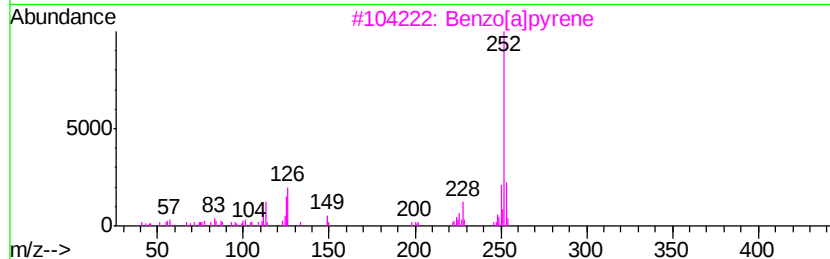
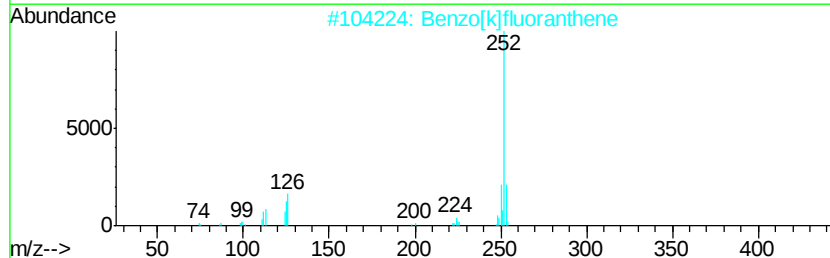
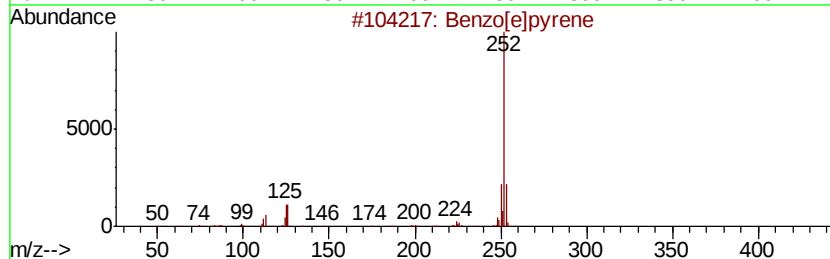
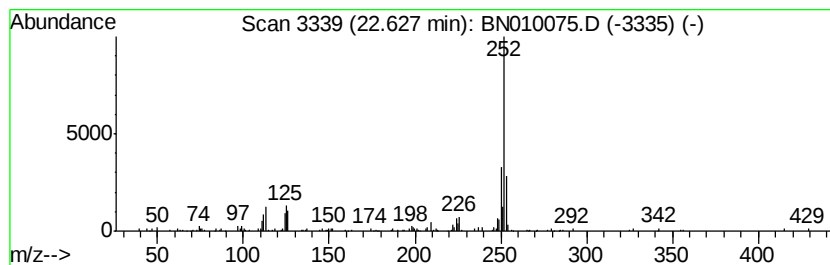
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzo[e]pyrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.63	3.34 ng/ul	202001	Perylene-d12	23.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	94
2		Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
3		Benzo[a]pyrene	252	C20H12	000050-32-8	89
4		Benzo[i]fluoranthene	252	C20H12	000205-82-3	89
5		Benzo[e]pyrene	252	C20H12	000192-97-2	86



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN030320\
Data File : BN010075.D
Acq On : 03 Mar 2020 23:22
Operator : CG/JU
Sample : L1619-04DL 10X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBDT8DL

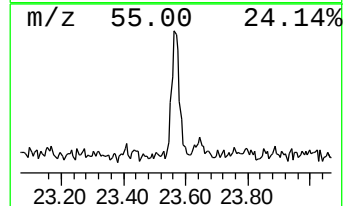
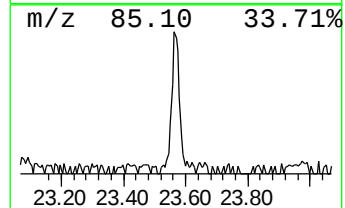
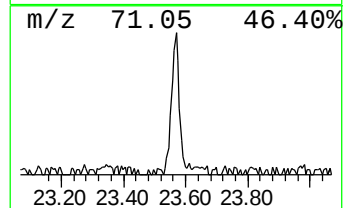
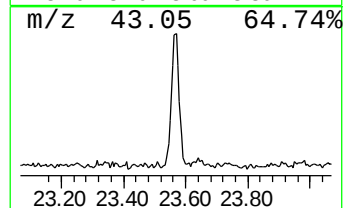
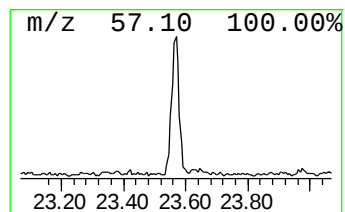
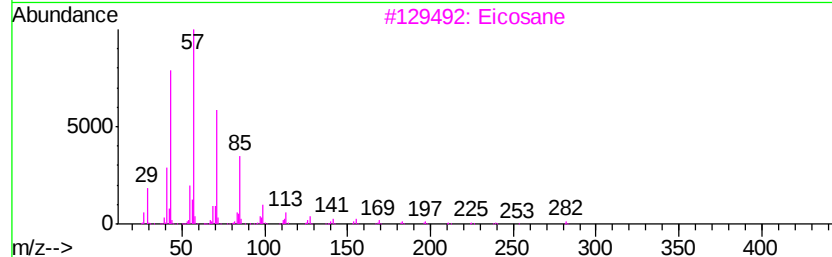
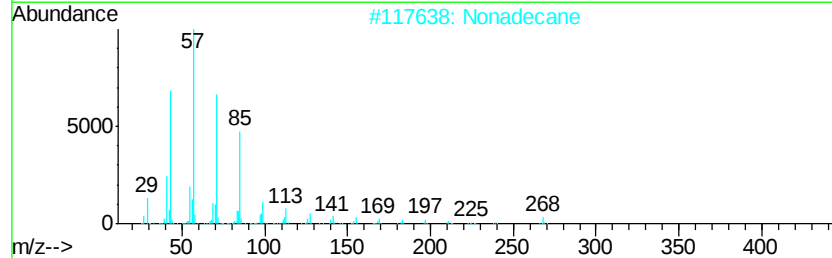
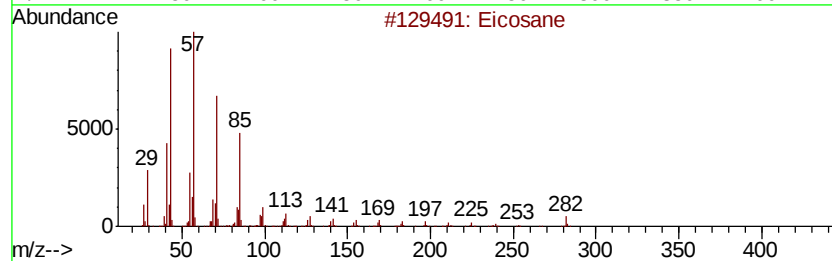
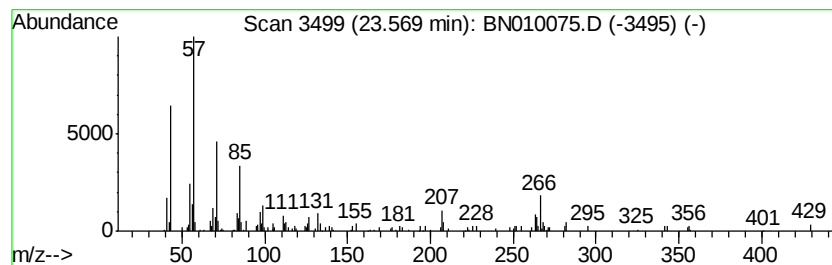
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.57	2.64 ng/ul	159590	Perylene-d12	23.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Eicosane	282	C20H42	000112-95-8	90
2		Nonadecane	268	C19H40	000629-92-5	78
3		Eicosane	282	C20H42	000112-95-8	78
4		2-methyloctacosane	408	C29H60	1000376-72-8	49
5		Tetratetracontane	619	C44H90	007098-22-8	49



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN030320\
Data File : BN010075.D
Acq On : 03 Mar 2020 23:22
Operator : CG/JU
Sample : L1619-04DL 10X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBDT8DL

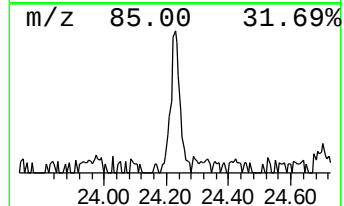
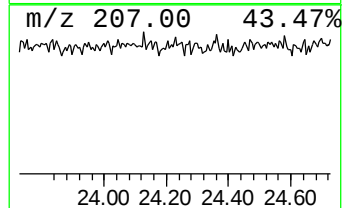
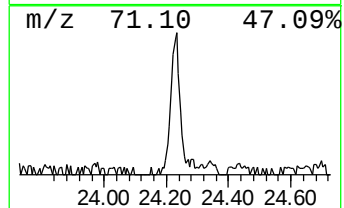
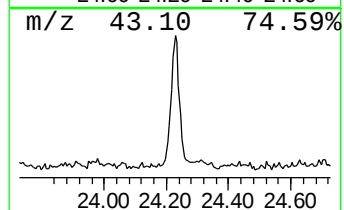
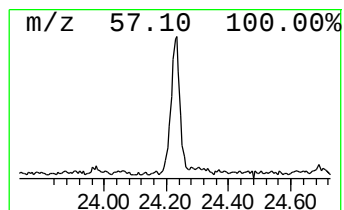
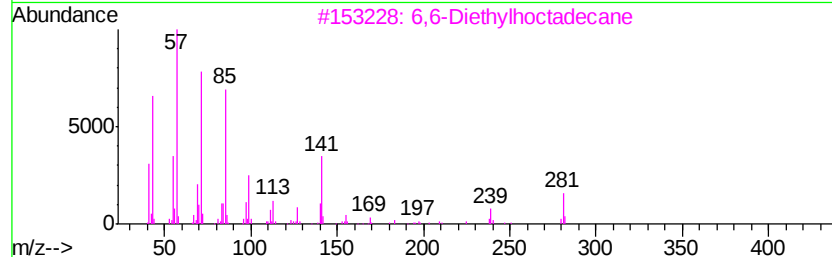
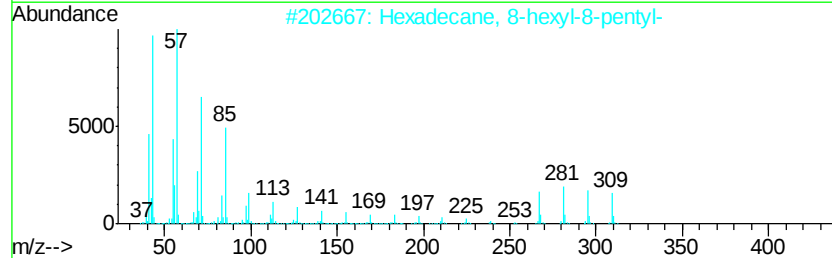
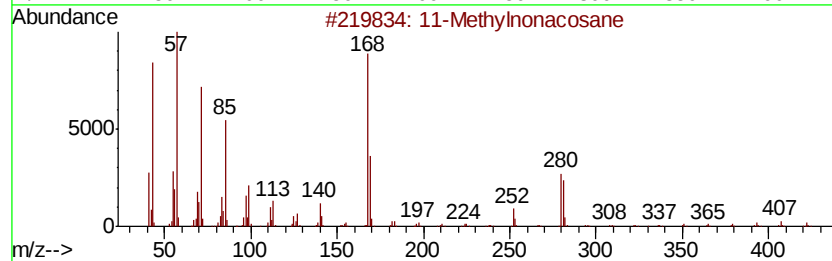
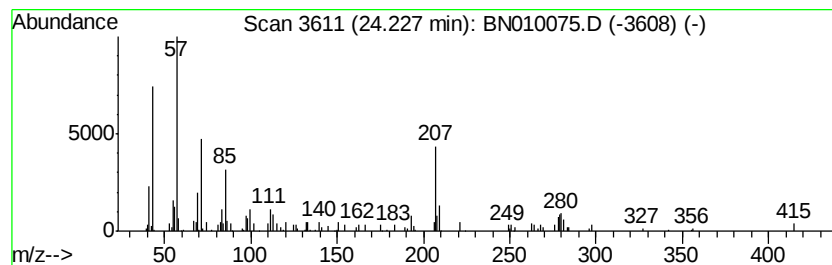
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.23	2.34 ng/ul	141335	Perylene-d12	23.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	11-Methylnonacosane	422	C30H62	007371-99-5	43
2		Hexadecane, 8-hexyl-8-pentyl-	380	C27H56	055282-29-6	40
3		6,6-Diethyloctadecane	310	C22H46	1000360-41-8	38
4		15-Methyltritriacontane	479	C34H70	1000131-20-3	37
5		Octadecane, 3-ethyl-5-(2-ethylbu...	366	C26H54	055282-12-7	35



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030320\
Data File : BN010075.D
Acq On : 03 Mar 2020 23:22
Operator : CG/JU
Sample : L1619-04DL 10X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBDT8DL

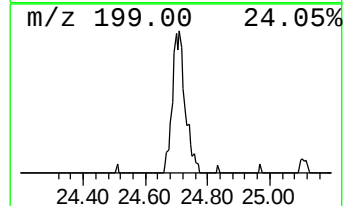
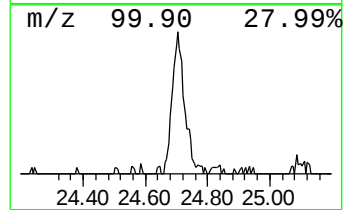
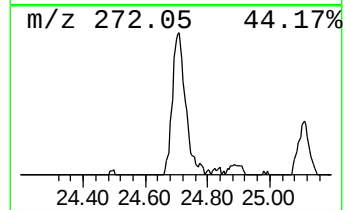
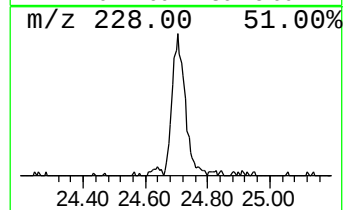
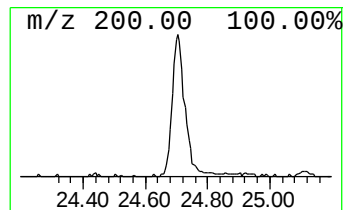
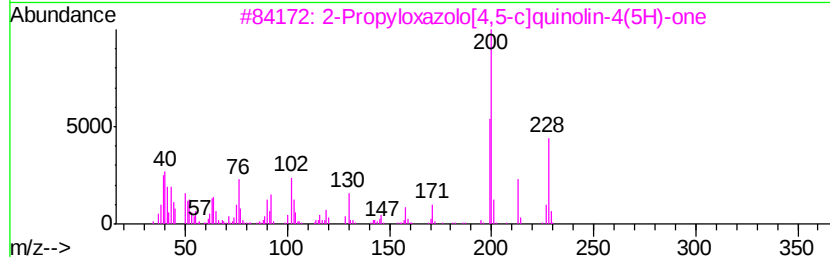
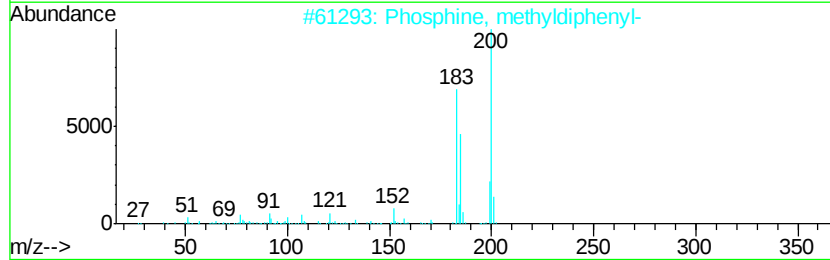
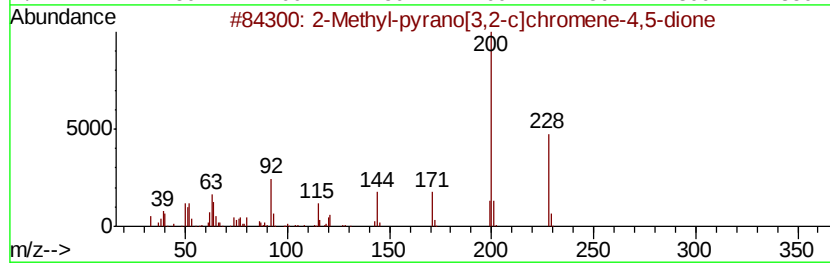
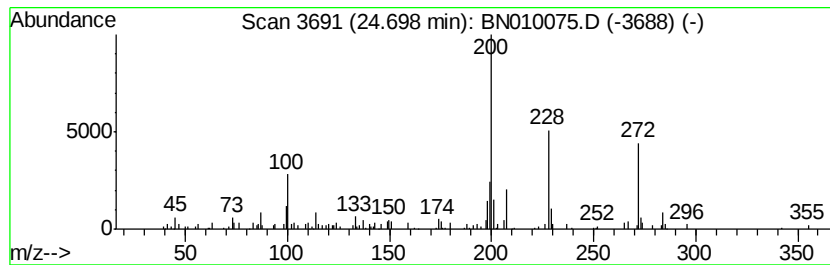
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.70	4.49 ng/ul	271570	Perylene-d12	23.07

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Methyl-pyrano[3,2-c]chromene-4...	228	C13H8O4	1000301-16-7	38
2		Phosphine, methylidiphenyl-	200	C13H13P	001486-28-8	27
3		2-Propyloxazolo[4,5-c]quinolin-4...	228	C13H12N2O2	211874-68-9	25
4		Ethylthiopentafluorobenzene	228	C8H5F5S	033288-21-0	22
5		N-[2-Methyl-4-pyridyl]-4-aminoph...	200	C12H12N2O	1000252-23-9	16



Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN030320\
Data File : BN010075.D
Acq On : 03 Mar 2020 23:22
Operator : CG/JU
Sample : L1619-04DL 10X
Misc :
ALS Vial : 19 Sample Multiplier: 1

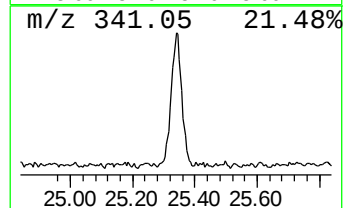
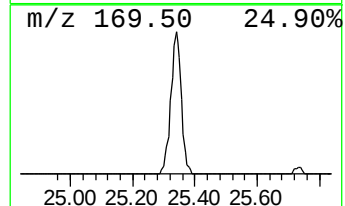
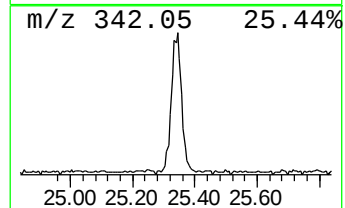
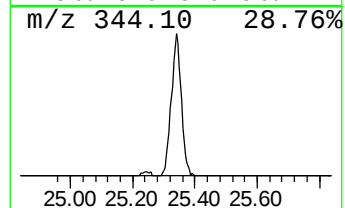
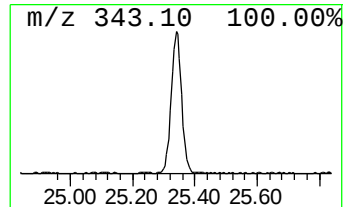
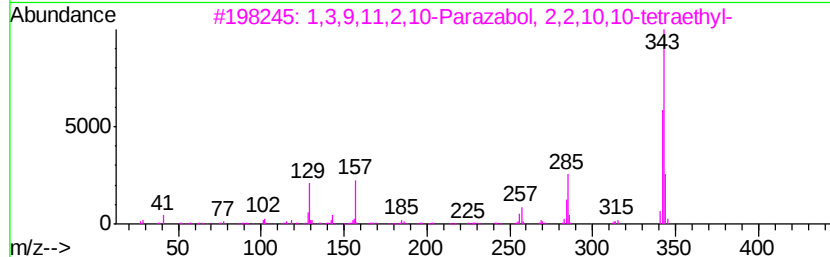
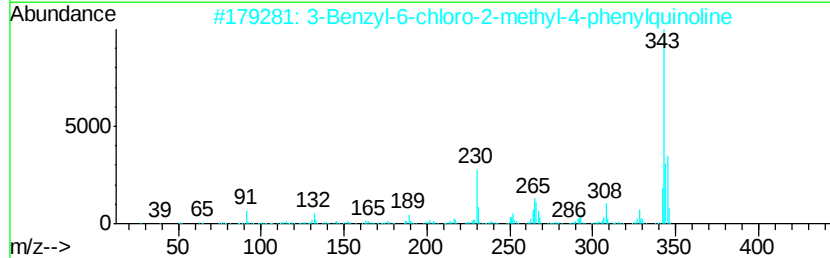
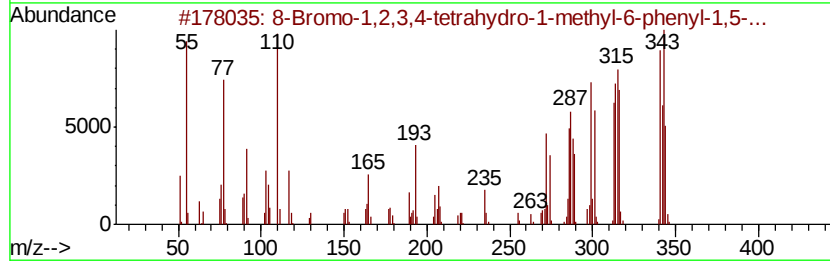
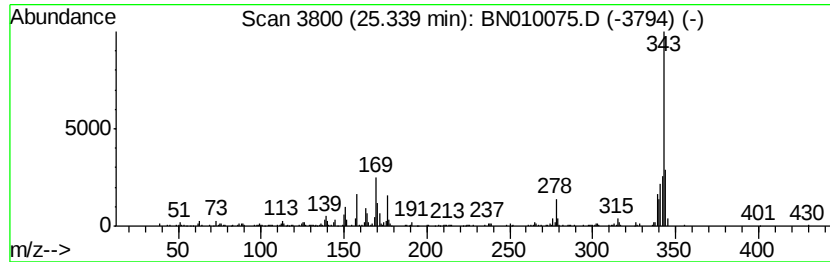
Instrument :
BNA_N
ClientSampleId :
DBDT8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 8-Bromo-1,2,3,4-tetrahydro-... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.		
25.34	13.66 ng/ul	826810	Perylene-d12		23.07		
Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	8-Bromo-1,2,3,4-tetrahydro-1-met...	342	C17H15BrN2O	063563-58-6	96		
2	3-Benzyl-6-chloro-2-methyl-4-phe...	343	C23H18ClN	022608-97-5	38		
3	1,3,9,11,2,10-Parazabol, 2,2,10,...	372	C22H30B2N4	1000157-80-0	37		
4	4-Cycloheptylaminomethylene-2-(4...	343	C20H26FN3O	1000276-03-9	35		
5	5-(p-Aminophenyl)-4-(4-biphenyly...	343	C21H17N3S	094863-57-7	32		



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN030320\
Data File : BN010075.D
Acq On : 03 Mar 2020 23:22
Operator : CG/JU
Sample : L1619-04DL 10X
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
DBDT8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN021220MA.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
Pyrene, 1-methyl-	19.69	2.2	ng/ul	226501	5	20.99	2067640	20.0
Cyclopenta[cd]pyrene	20.70	2.2	ng/ul	224575	5	20.99	2067640	20.0
Anthracene, 9-phe...	22.24	4.9	ng/ul	294351	6	23.07	1210220	20.0
Benzo[e]pyrene	22.63	3.3	ng/ul	202001	6	23.07	1210220	20.0
(DEL) Alkane: Str...	23.57	2.6	ng/ul	159590	6	23.07	1210220	20.0
(DEL) Alkane: Str...	24.23	2.3	ng/ul	141335	6	23.07	1210220	20.0
unknown-01	24.70	4.5	ng/ul	271570	6	23.07	1210220	20.0
8-Bromo-1,2,3,4-t...	25.34	13.7	ng/ul	826810	6	23.07	1210220	20.0