

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN100618\
 Data File : BN002974.D
 Acq On : 05 Oct 2018 21:55
 Operator : MJ/SJ
 Sample : J5230-04
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A3YY0

Integration Parameters: LSCINT.P

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

Filtering: 5
 Min Area: 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-BN091418MA.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	6.822	646	652	667	rVB	332450	628111	14.92%	1.150%
2	6.969	670	677	691	rBV	287034	462681	10.99%	0.847%
3	7.169	704	711	721	rBV	357547	605005	14.37%	1.108%
4	7.622	780	788	796	rBV	456305	748107	17.77%	1.370%
5	8.345	904	911	923	rBV	407991	738234	17.54%	1.352%
6	8.775	977	984	992	rBV	352864	602523	14.31%	1.104%
7	9.492	1099	1106	1116	rBV	308086	540731	12.84%	0.990%
8	10.040	1192	1199	1215	rBV	435269	840521	19.97%	1.539%
9	10.392	1250	1259	1266	rBV	713281	1233473	29.30%	2.259%
10	10.545	1278	1285	1301	rBV	335419	648046	15.39%	1.187%
11	13.086	1712	1717	1723	rBV2	33589	50482	1.20%	0.092%
12	13.163	1723	1730	1738	rVV	204188	333264	7.92%	0.610%
13	13.681	1812	1818	1822	rBV	868764	1267690	30.11%	2.322%
14	13.728	1822	1826	1833	rVB	664709	961941	22.85%	1.762%
15	13.951	1857	1864	1879	rBV2	1072634	1991022	47.30%	3.647%
16	14.169	1895	1901	1905	rBV	51905	69839	1.66%	0.128%
17	14.263	1910	1917	1924	rBV2	1113854	1691038	40.17%	3.097%
18	14.522	1952	1961	1976	rBV	215508	557072	13.23%	1.020%
19	15.128	2060	2064	2071	rVB2	44757	69780	1.66%	0.128%
20	15.263	2081	2087	2093	rBV	1419168	2027603	48.16%	3.714%
21	15.316	2093	2096	2106	rVB5	27897	60466	1.44%	0.111%
22	15.416	2107	2113	2122	rBV	375335	610053	14.49%	1.117%
23	16.004	2206	2213	2219	rVB5	61457	150043	3.56%	0.275%
24	16.563	2304	2308	2311	rBV3	26431	44700	1.06%	0.082%
25	16.680	2323	2328	2334	rBV2	44772	69334	1.65%	0.127%
26	16.780	2342	2345	2351	rVV2	31129	59017	1.40%	0.108%
27	16.839	2351	2355	2359	rVB2	32363	44921	1.07%	0.082%
28	17.016	2379	2385	2390	rBV	1376384	2079091	49.39%	3.808%
29	17.057	2390	2392	2397	rVV	396042	525224	12.48%	0.962%
30	17.116	2397	2402	2406	rVV2	1595339	2257463	53.62%	4.135%
31	17.151	2406	2408	2414	rVV	211678	279095	6.63%	0.511%
32	17.210	2414	2418	2424	rVV5	28808	55098	1.31%	0.101%
33	17.433	2451	2456	2460	rBV	55213	105019	2.49%	0.192%
34	17.521	2468	2471	2482	rVB2	36963	81460	1.94%	0.149%

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35	17.610	2482	2486	2488	rBV4	24256	43392	1.03%	0.079%
36	17.639	2488	2491	2497	rVB	65421	80865	1.92%	0.148%
37	17.898	2527	2535	2536	rBV	96731	140918	3.35%	0.258%
38	17.927	2536	2540	2542	rBV	153916	217835	5.17%	0.399%
39	17.992	2547	2551	2554	rBV	53936	77808	1.85%	0.143%
40	18.021	2554	2556	2561	rVB	55132	68374	1.62%	0.125%
41	18.086	2562	2567	2571	rBV2	140999	245442	5.83%	0.450%
42	18.216	2584	2589	2593	rBV4	55981	89597	2.13%	0.164%
43	18.398	2616	2620	2625	rBV	94758	144366	3.43%	0.264%
44	18.451	2625	2629	2638	rVB	94405	142784	3.39%	0.262%
45	18.698	2668	2671	2675	rVB	46724	61799	1.47%	0.113%
46	18.739	2675	2678	2682	rVB	53124	62310	1.48%	0.114%
47	18.821	2682	2692	2695	rBV2	138491	272967	6.48%	0.500%
48	18.857	2695	2698	2705	rVV5	74912	192182	4.57%	0.352%
49	18.916	2705	2708	2711	rVB	50208	58883	1.40%	0.108%
50	18.968	2712	2717	2723	rBV2	149844	256551	6.09%	0.470%
51	19.086	2731	2737	2744	rVB	2187754	2890093	68.65%	5.293%
52	19.151	2745	2748	2751	rBV	56176	67657	1.61%	0.124%
53	19.186	2752	2754	2757	rVV4	46520	60085	1.43%	0.110%
54	19.239	2760	2763	2767	rVB	120194	145777	3.46%	0.267%
55	19.421	2789	2794	2797	rBV2	2041202	3190616	75.79%	5.844%
56	19.451	2797	2799	2808	rVV	1905993	2332361	55.40%	4.272%
57	19.527	2809	2812	2820	rVB3	113566	227159	5.40%	0.416%
58	19.633	2826	2830	2837	rBV	108911	186856	4.44%	0.342%
59	19.780	2850	2855	2863	rBV4	176466	427465	10.15%	0.783%
60	19.915	2874	2878	2881	rBV4	148804	271317	6.44%	0.497%
61	20.051	2899	2901	2904	rVB2	83483	82570	1.96%	0.151%
62	20.110	2906	2911	2915	rVV2	205995	348204	8.27%	0.638%
63	20.251	2932	2935	2938	rBV	104757	137633	3.27%	0.252%
64	20.710	3009	3013	3018	rVB	390600	519301	12.34%	0.951%
65	20.868	3036	3040	3043	rBV2	279242	374425	8.89%	0.686%
66	20.904	3043	3046	3049	rVV	219736	264146	6.27%	0.484%
67	20.945	3049	3053	3059	rVV3	530689	928291	22.05%	1.700%
68	21.004	3060	3063	3070	rVB	295567	417796	9.92%	0.765%
69	21.221	3094	3100	3104	rBV2	2188745	4209784	100.00%	7.711%
70	21.262	3104	3107	3116	rVB	1232948	1620474	38.49%	2.968%
71	21.380	3124	3127	3129	rBV2	142086	184079	4.37%	0.337%

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72	21.810	3197	3200	3202	rBV	158440	219260	5.21%	0.402%
73	22.257	3273	3276	3281	rVB2	402823	501668	11.92%	0.919%
74	22.780	3357	3365	3370	rBV2	1131588	2835433	67.35%	5.193%
75	22.945	3390	3393	3399	rVB	226637	341175	8.10%	0.625%
76	23.245	3440	3444	3448	rBV	451651	763871	18.15%	1.399%
77	23.298	3448	3453	3458	rVV2	1351420	2605130	61.88%	4.771%
78	23.339	3458	3460	3468	rVB	624215	981239	23.31%	1.797%
79	23.439	3471	3477	3482	rBV	890823	1613074	38.32%	2.954%
80	23.927	3556	3560	3566	rVB	358975	618921	14.70%	1.134%
81	24.651	3679	3683	3693	rVB	295585	587871	13.96%	1.077%

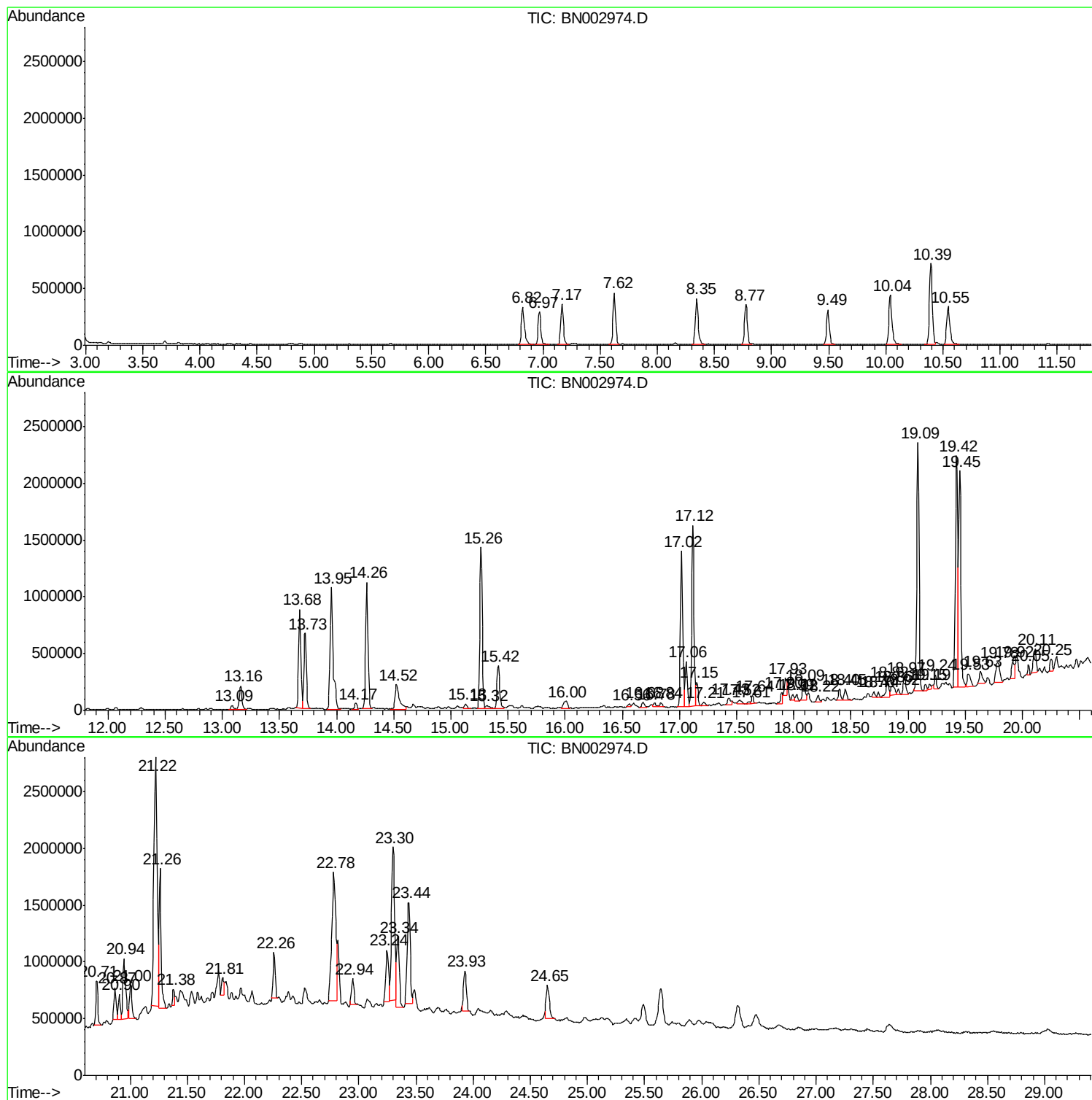
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091418MA.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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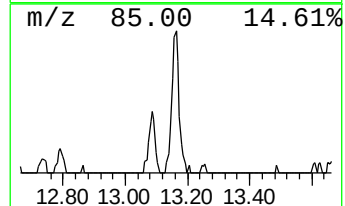
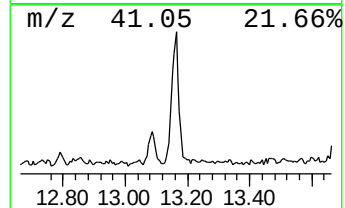
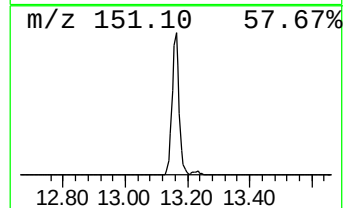
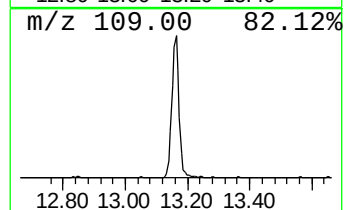
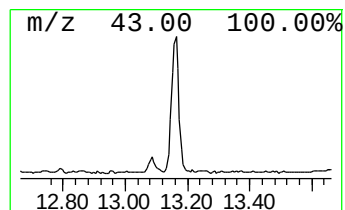
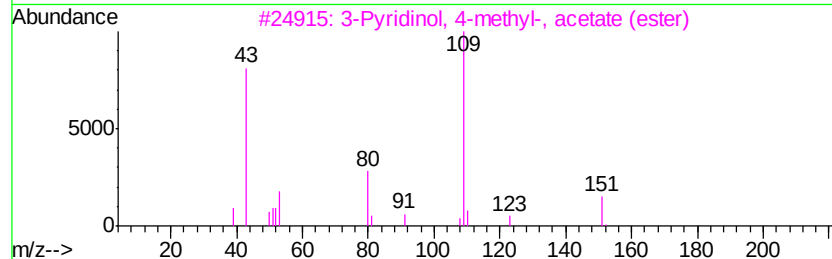
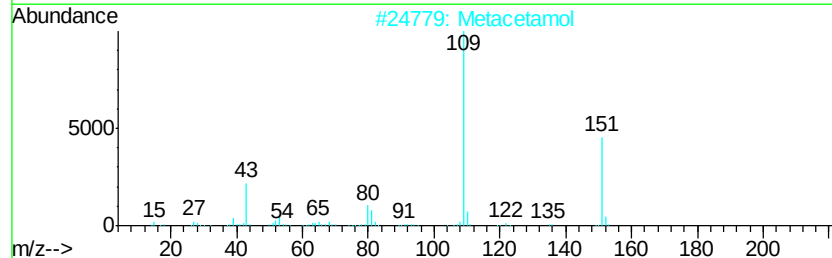
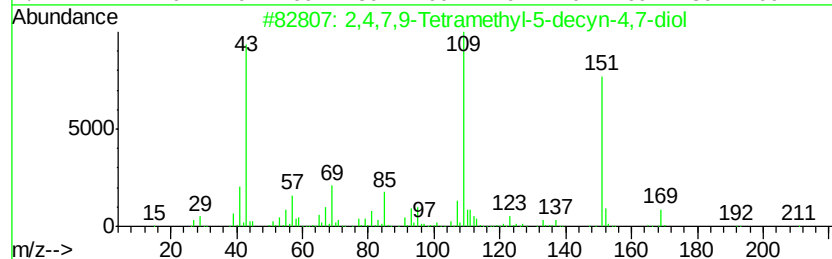
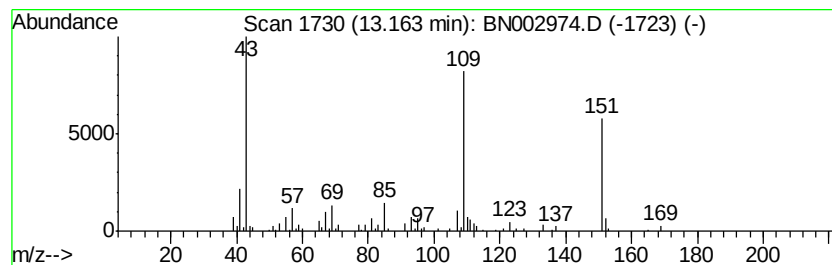
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091418MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.16	3.94 ng/ul	333264	Acenaphthene-d10	14.26

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	91		
2	Metacetamol	151	C8H9NO2	000621-42-1	59		
3	3-Pyridinol, 4-methyl-, acetate ...	151	C8H9NO2	001006-96-8	59		
4	Pyridine, 3-butyl-, 1-oxide	151	C9H13NO	031396-33-5	59		
5	Cyclohexanol, 3,3,5-trimethyl-	142	C9H18O	000116-02-9	43		



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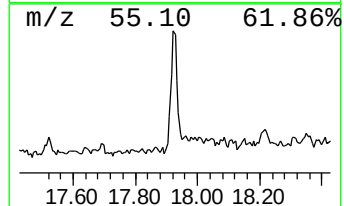
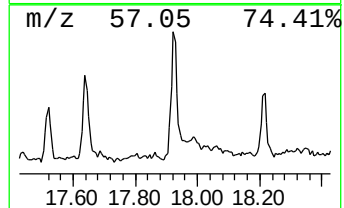
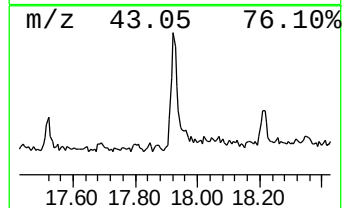
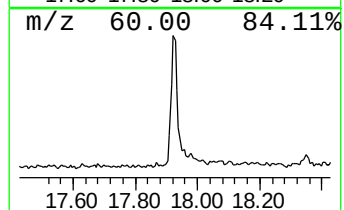
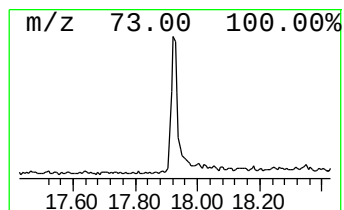
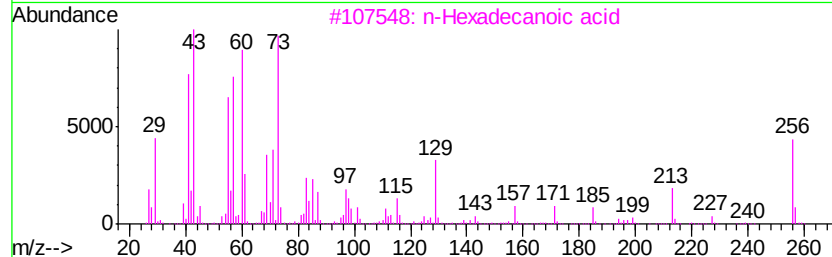
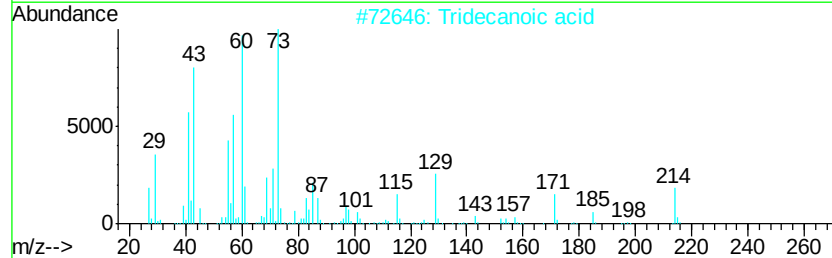
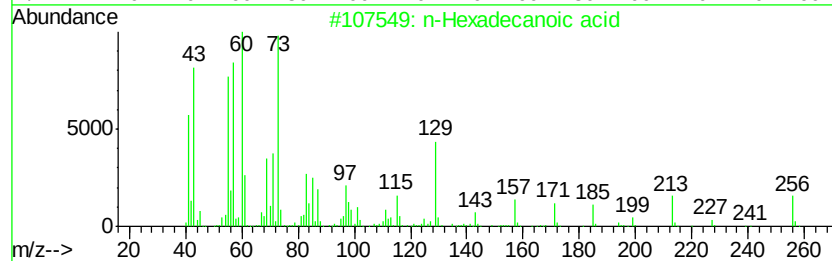
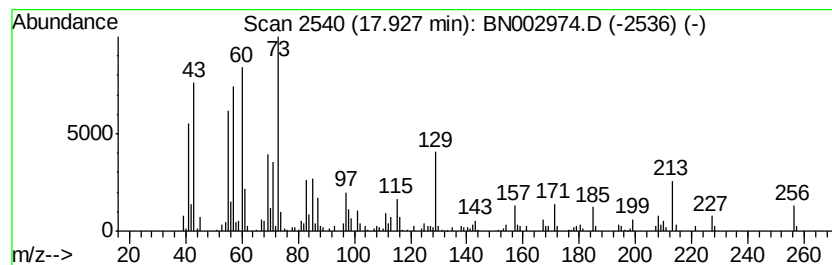
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 n-Hexadecanoic acid Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.93	2.10 ng/ul	217835	Phenanthrene-d10		17.02	
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	99
2	Tridecanoic acid		214	C13H26O2	000638-53-9	93
3	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	91
4	Tetradecanoic acid		228	C14H28O2	000544-63-8	89
5	Tetradecanoic acid		228	C14H28O2	000544-63-8	81



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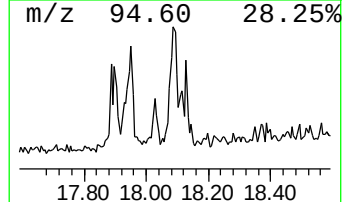
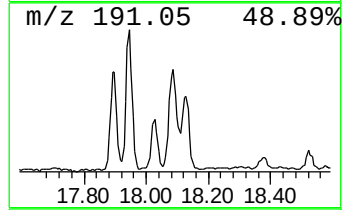
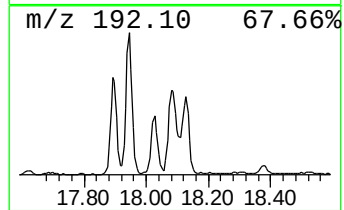
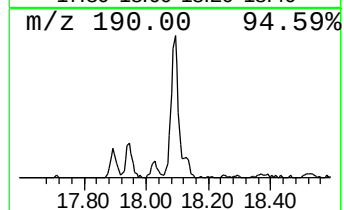
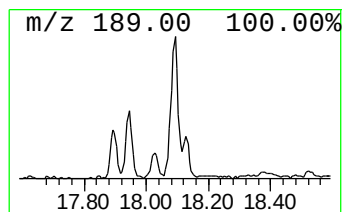
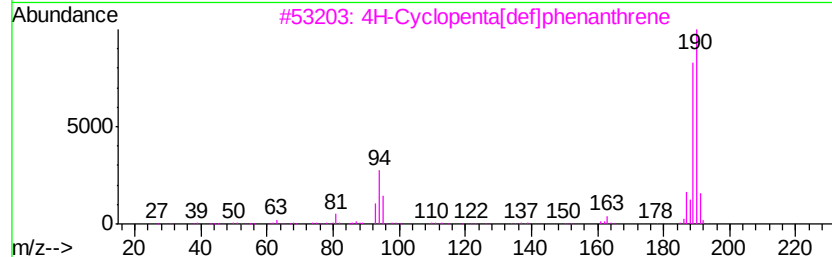
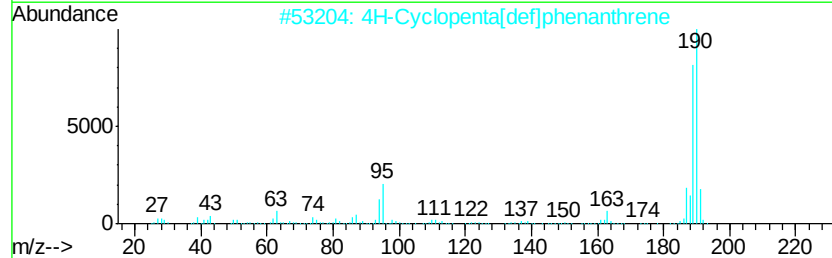
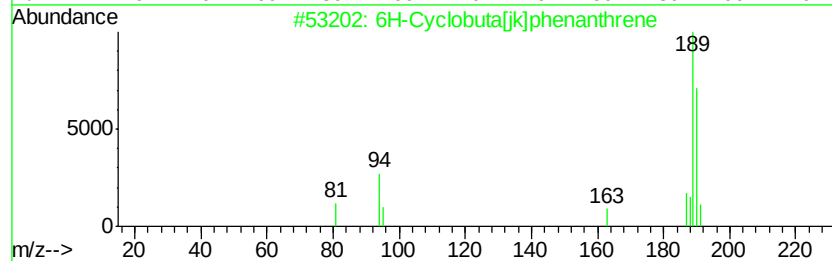
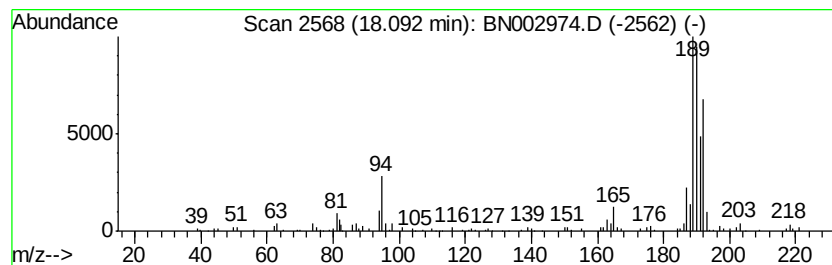
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Peak Number 3 unknown-01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.09	2.36 ng/ul	245442	Phenanthrene-d10	17.02	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	46
2	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	43
3	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	43
4	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	42
5	1a,9b-Dihydro-1H-cyclopropa[a]an...	192	C15H12	006109-24-6	42



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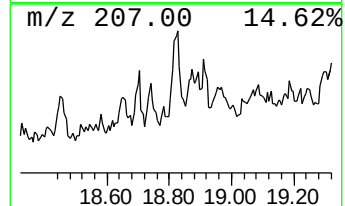
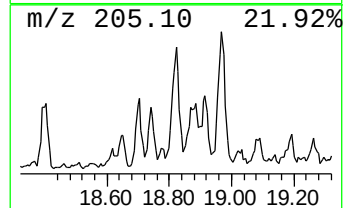
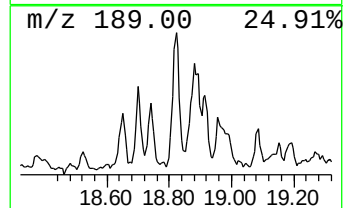
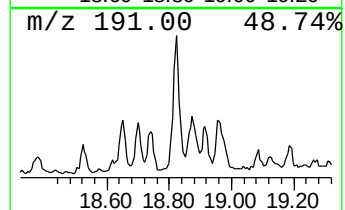
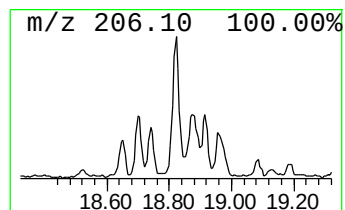
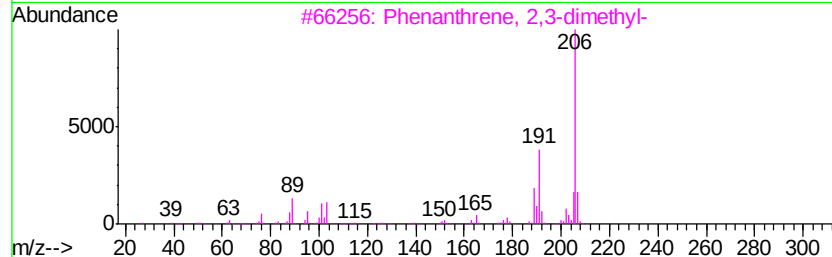
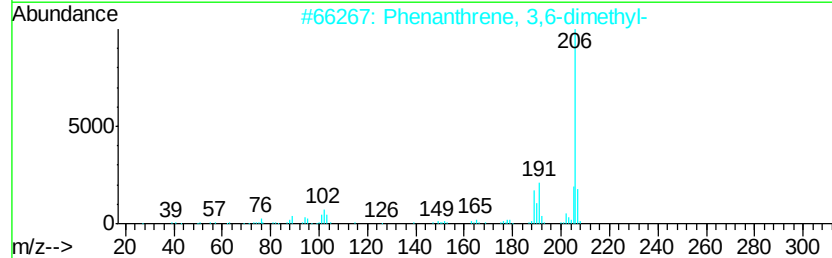
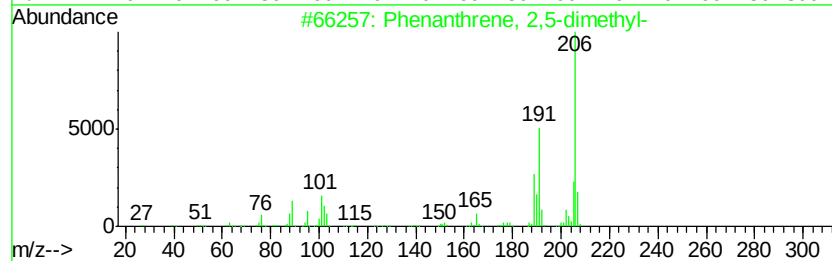
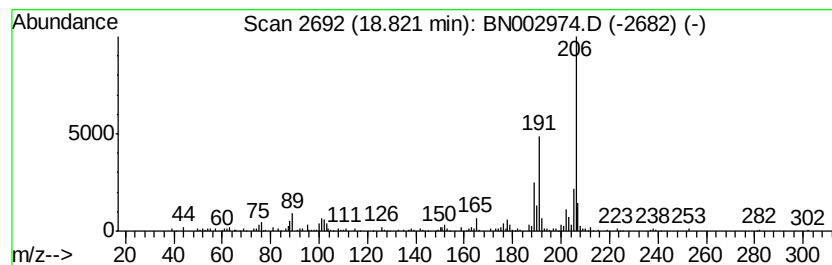
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Phenanthrene, 2,5-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.82	2.63 ng/ul	272967	Phenanthrene-d10	17.02		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	96	
2	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	90	
3	Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	87	
4	Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	86	
5	Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	81	



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Data File : BN002974.D
Acq On : 05 Oct 2018 21:55
Operator : MJ/SJ
Sample : J5230-04
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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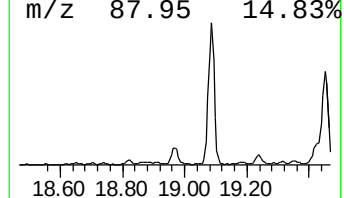
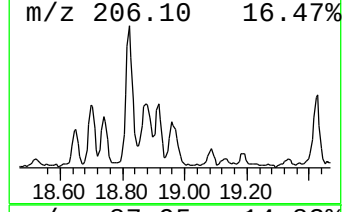
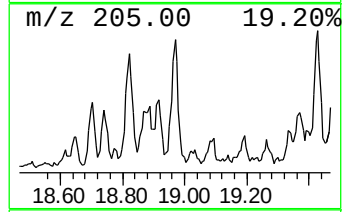
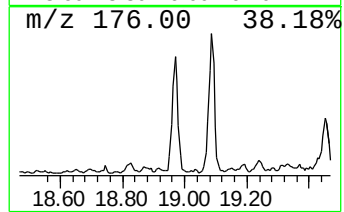
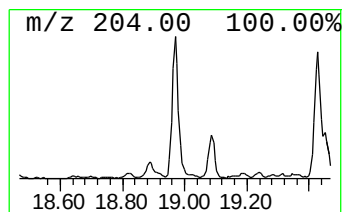
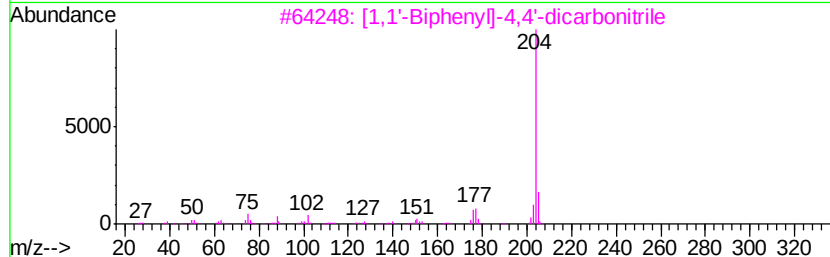
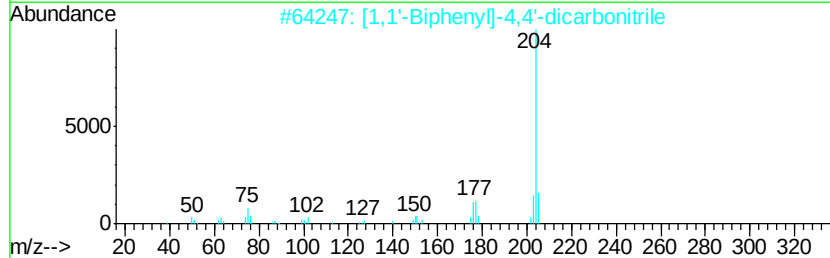
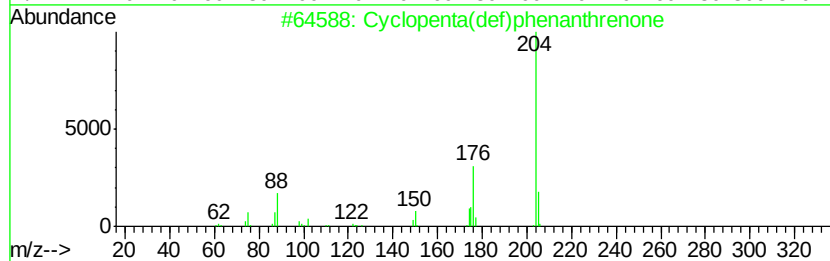
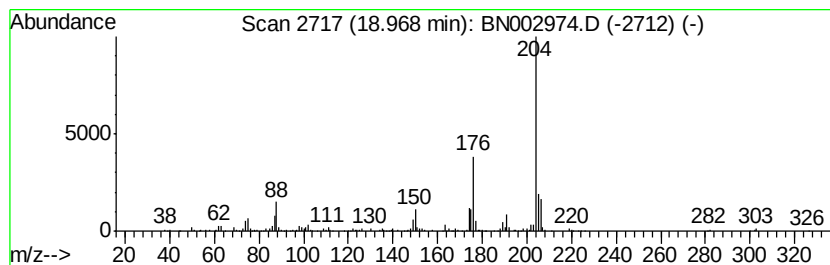
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Cyclopenta(def)phenanthrenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.97	2.47 ng/ul	256551	Phenanthrene-d10	17.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopenta(def)phenanthrenone	204	C15H8O	005737-13-3	96
2		[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	50
3		[1,1'-Biphenyl]-4,4'-dicarbonitrile	204	C14H8N2	001591-30-6	49
4		[1,1'-Biphenyl]-2,2'-dicarbonitrile	204	C14H8N2	004341-02-0	49
5		Naphthalene, 2-phenyl-	204	C16H12	000612-94-2	47



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Sample : J5230-04
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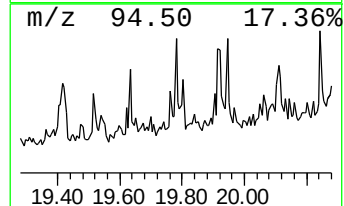
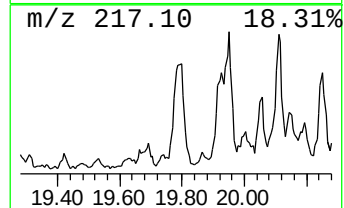
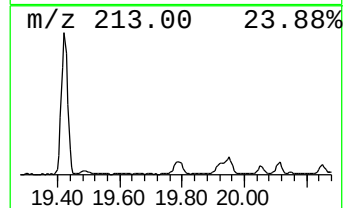
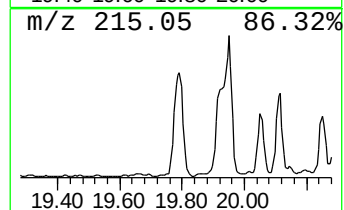
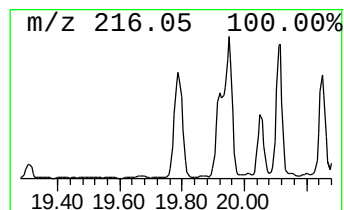
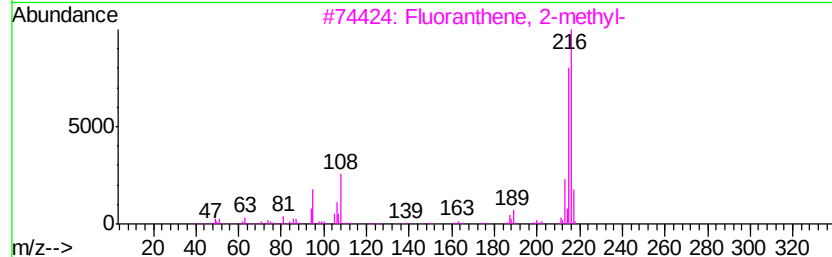
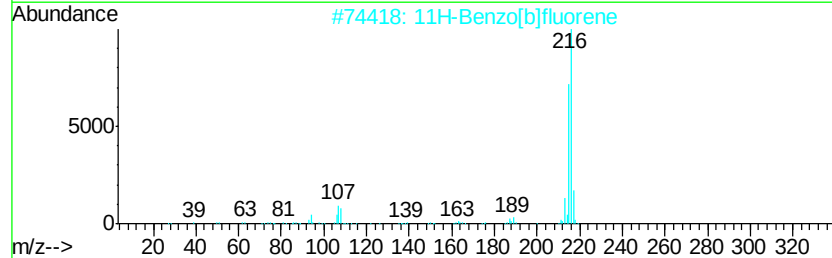
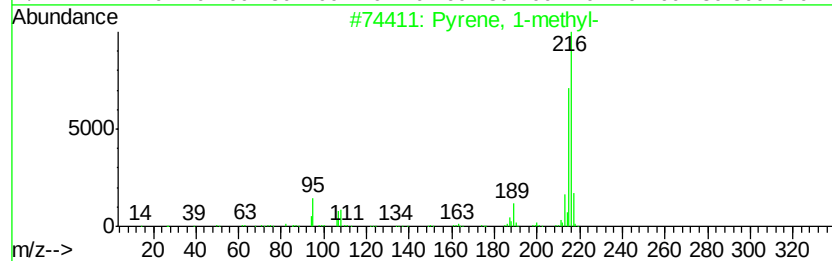
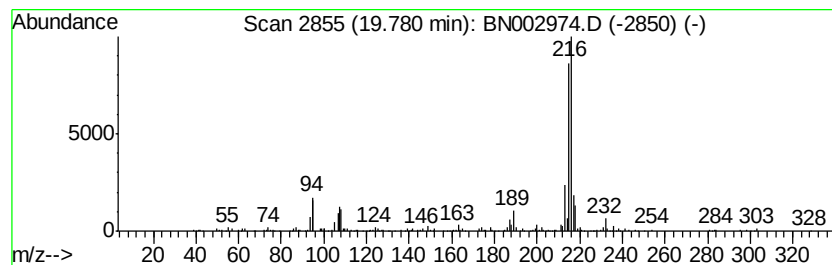
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Pyrene, 1-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.78	2.03 ng/ul	427465	Chrysene-d12	21.23
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Pyrene, 1-methyl-	216 C17H12	002381-21-7 95
2		11H-Benzo[b]fluorene	216 C17H12	000243-17-4 90
3		Fluoranthene, 2-methyl-	216 C17H12	033543-31-6 90
4		Pyrene, 1-methyl-	216 C17H12	002381-21-7 89
5		Pyrene, 1-methyl-	216 C17H12	002381-21-7 89



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN100618\
Data File : BN002974.D
Acq On : 05 Oct 2018 21:55
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Sample : J5230-04
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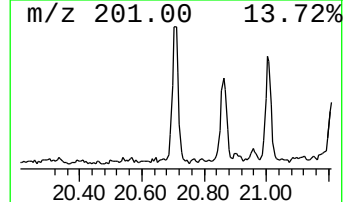
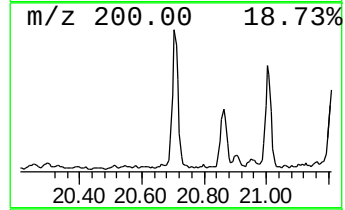
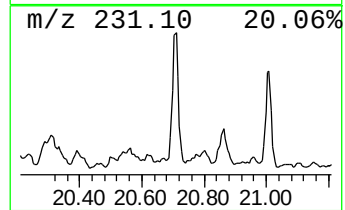
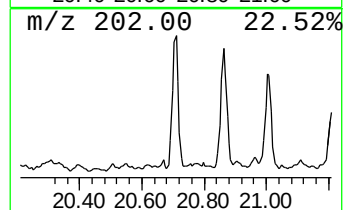
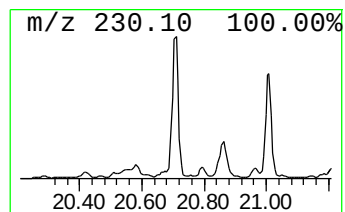
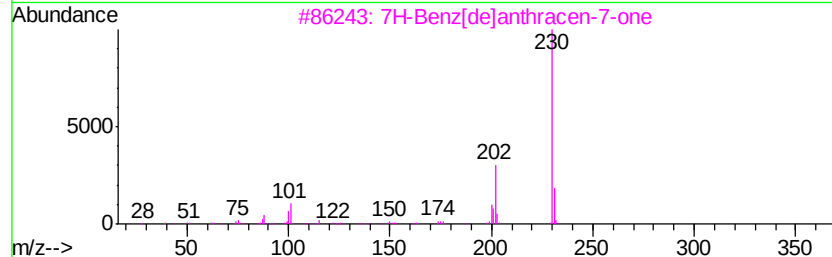
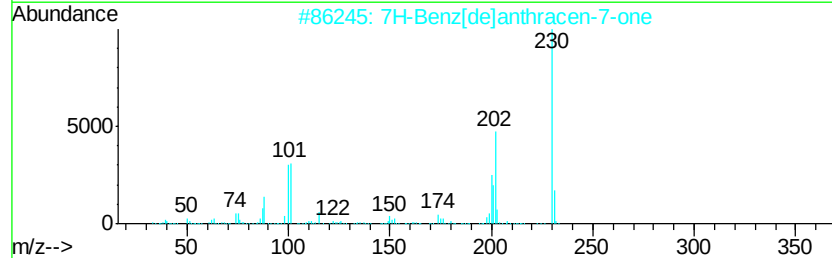
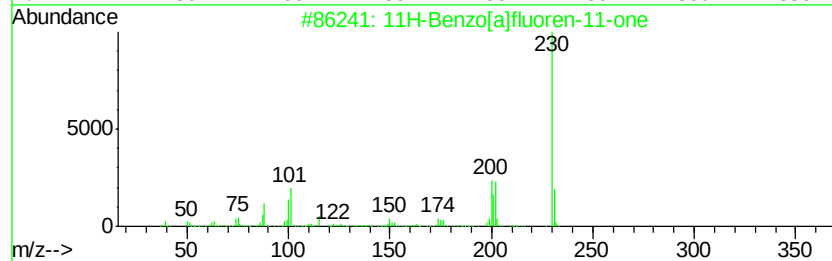
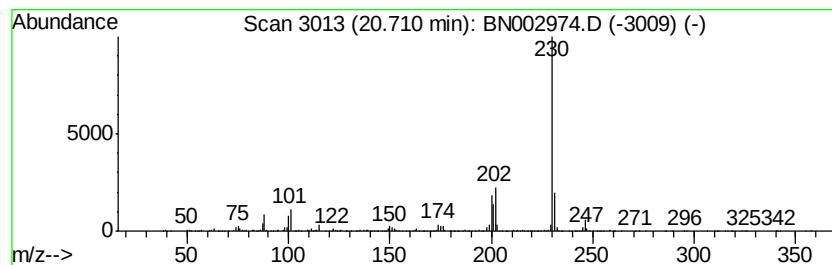
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 11H-Benzo[a]fluoren-11-one Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.71	2.47 ng/ul	519301	Chrysene-d12	21.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	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	90
4		7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	90
5		7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	76



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN100618\
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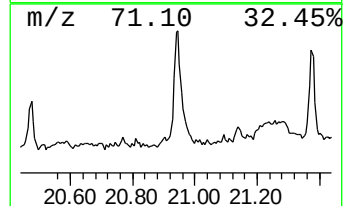
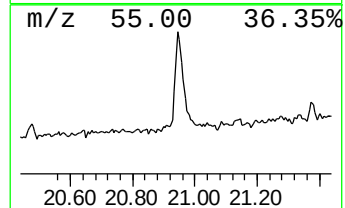
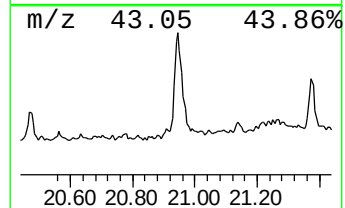
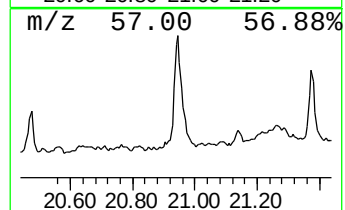
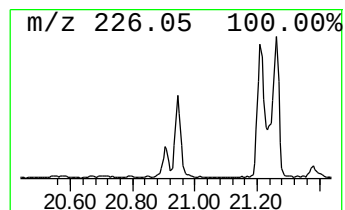
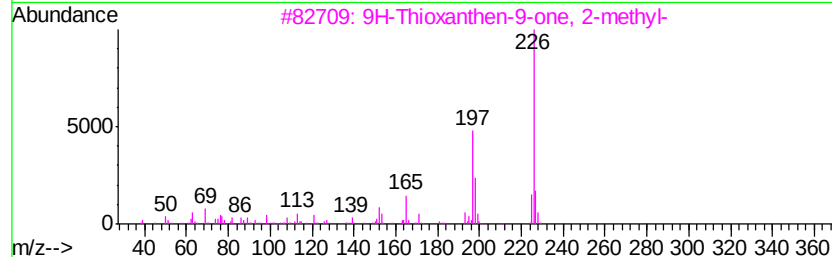
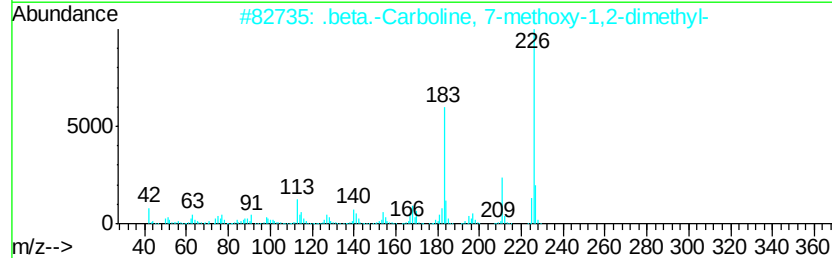
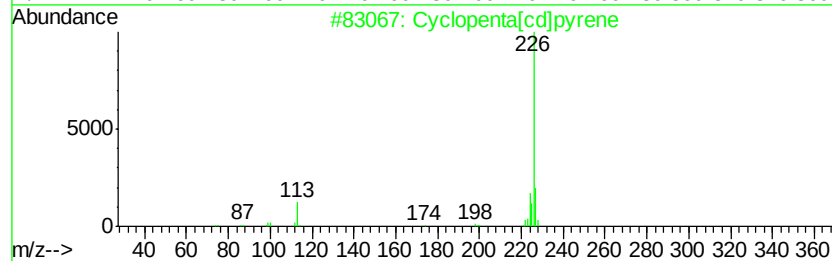
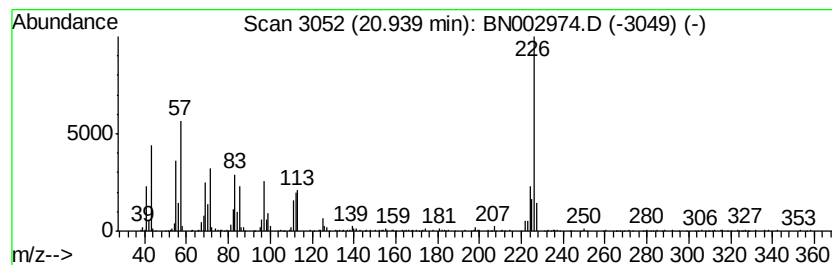
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-02 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.94	4.41 ng/ul	928291	Chrysene-d12	21.23
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Cyclopenta[cd]pyrene	226 C18H10	027208-37-3 49
2		.beta.-Carboline, 7-methoxy-1,2-...	226 C14H14N2O	006519-18-2 49
3		9H-Thioxanthen-9-one, 2-methyl-	226 C14H10OS	015774-82-0 47
4		Isonipecotic acid, N-(cyclohexyl...	337 C20H35NO3	1000361-03-0 43
5		2-Oxa-7-azatricyclo[4.4.0.0(3,8)]...	381 C18H23NO6S	055905-56-1 40



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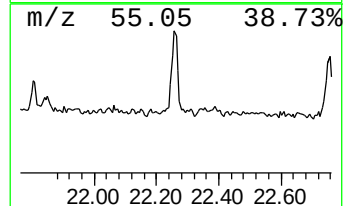
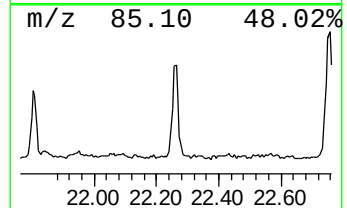
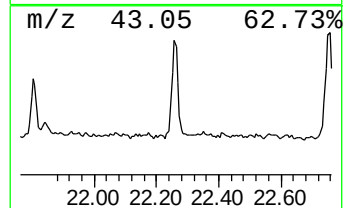
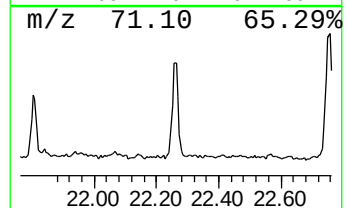
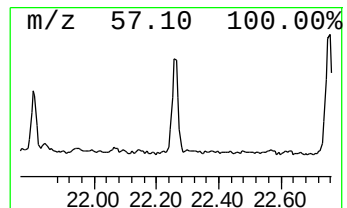
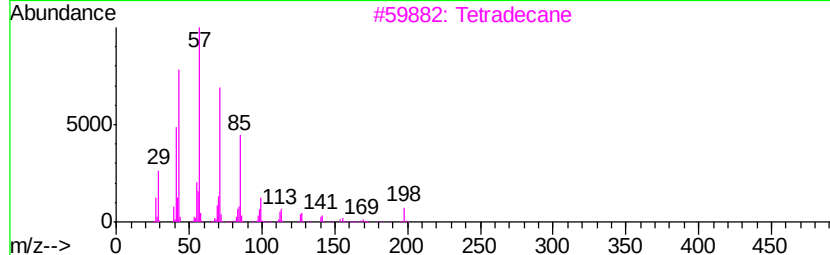
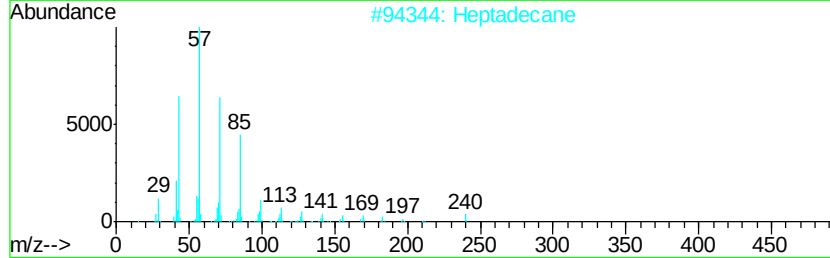
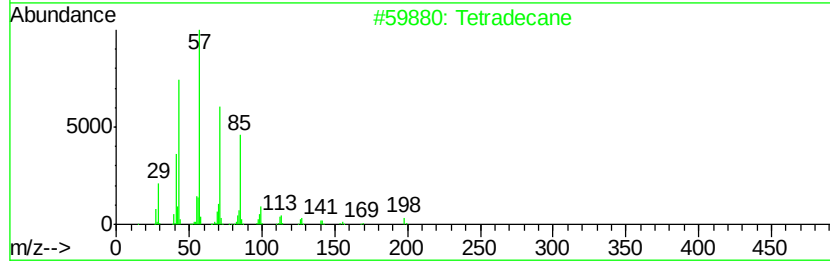
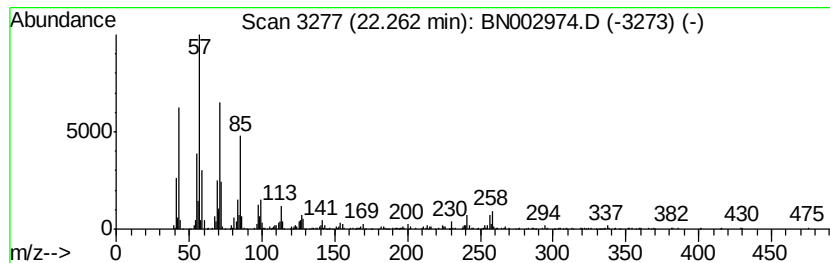
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.26	2.38 ng/ul	501668	Chrysene-d12	21.23

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetradecane	198	C14H30	000629-59-4	89
2		Heptadecane	240	C17H36	000629-78-7	87
3		Tetradecane	198	C14H30	000629-59-4	70
4		Heptadecane	240	C17H36	000629-78-7	70
5		Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	55



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN100618\
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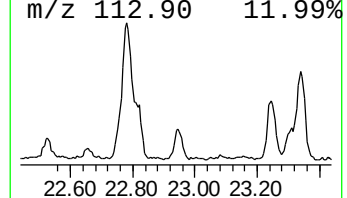
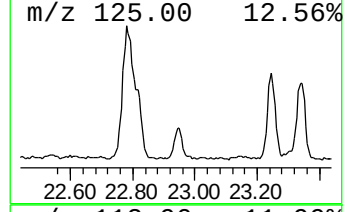
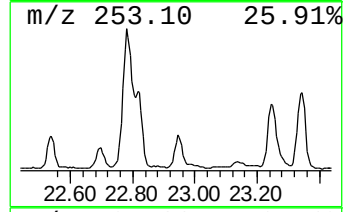
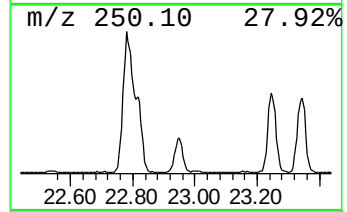
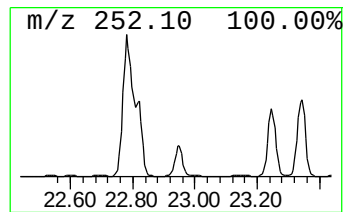
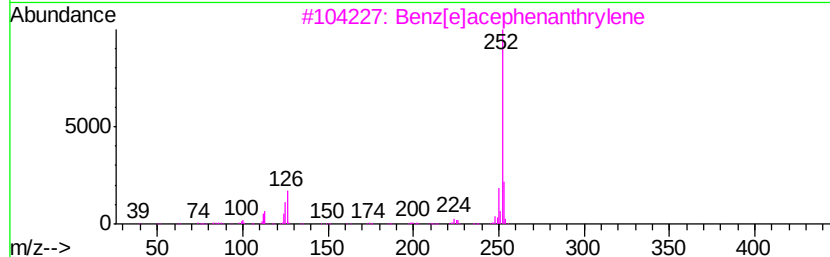
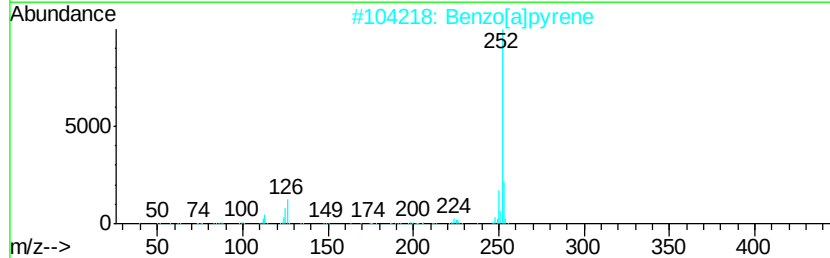
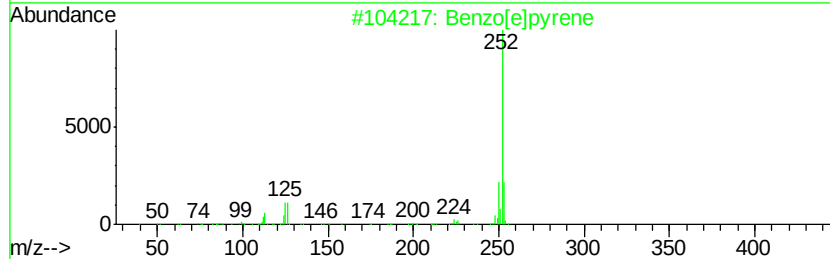
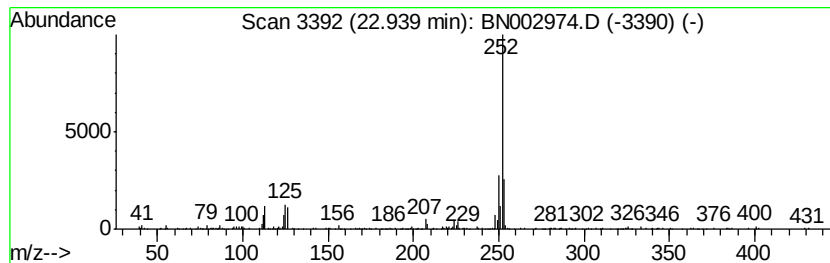
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Benzo[e]pyrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.94	4.23 ng/ul	341175	Perylene-d12	23.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzo[e]pyrene	252	C20H12	000192-97-2	96
2		Benzo[a]pyrene	252	C20H12	000050-32-8	93
3		Benz[e]acephenanthrylene	252	C20H12	000205-99-2	93
4		Perylene	252	C20H12	000198-55-0	93
5		Benzo[k]fluoranthene	252	C20H12	000207-08-9	90



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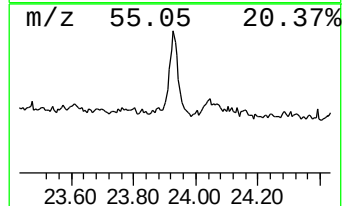
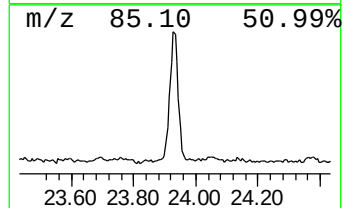
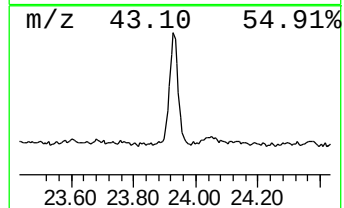
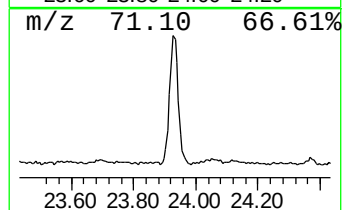
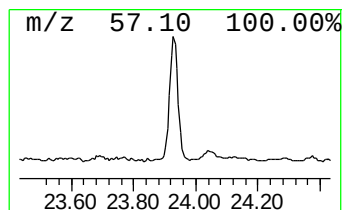
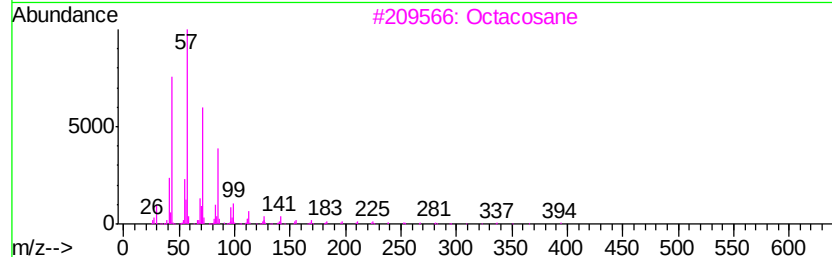
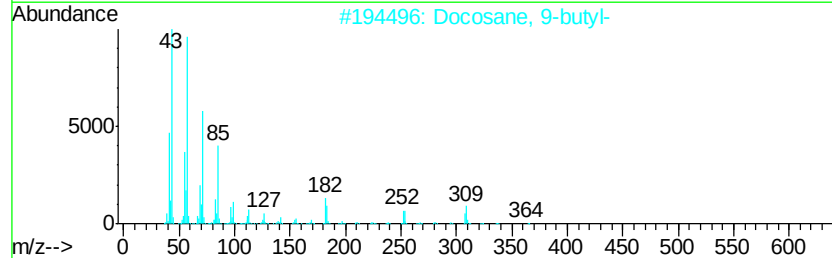
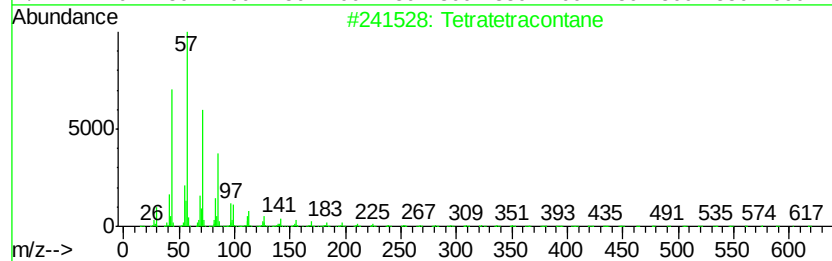
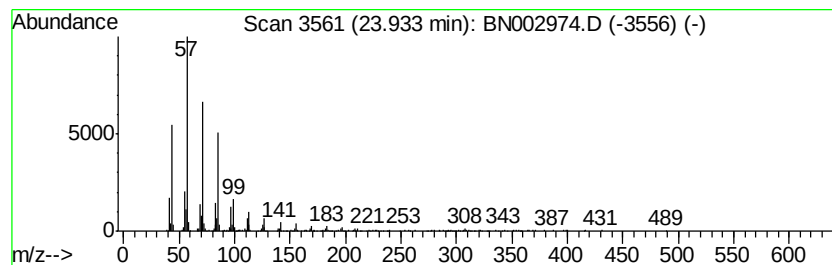
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.93	7.67 ng/ul	618921	Perylene-d12	23.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Tetratetracontane	619	C44H90	007098-22-8	91
2		Docosane, 9-butyl-	366	C26H54	055282-14-9	90
3		Octacosane	394	C28H58	000630-02-4	90
4		Heneicosane	296	C21H44	000629-94-7	90
5		Heptacosane	380	C27H56	000593-49-7	87



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN100618\
Data File : BN002974.D
Acq On : 05 Oct 2018 21:55
Operator : MJ/SJ
Sample : J5230-04
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
A3YY0

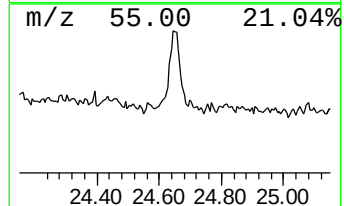
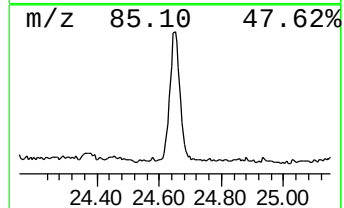
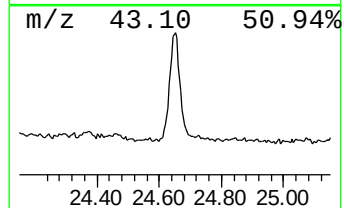
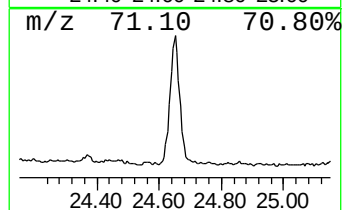
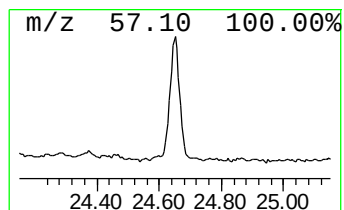
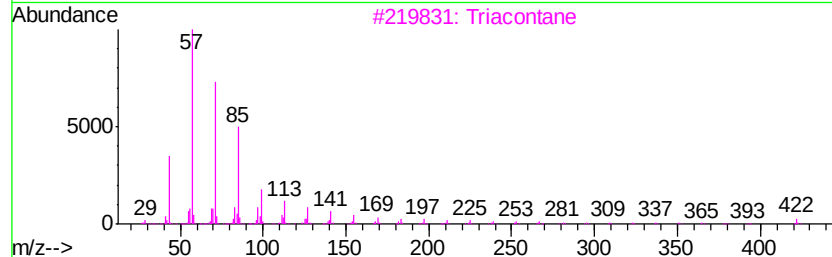
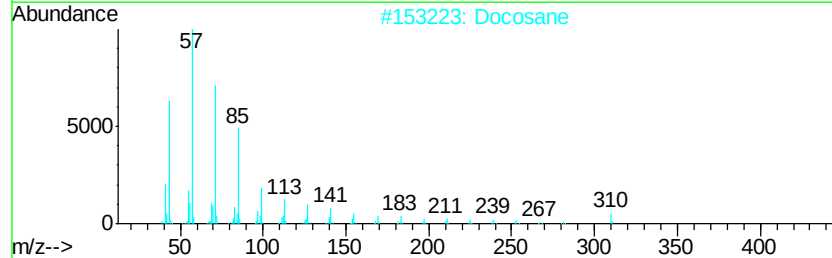
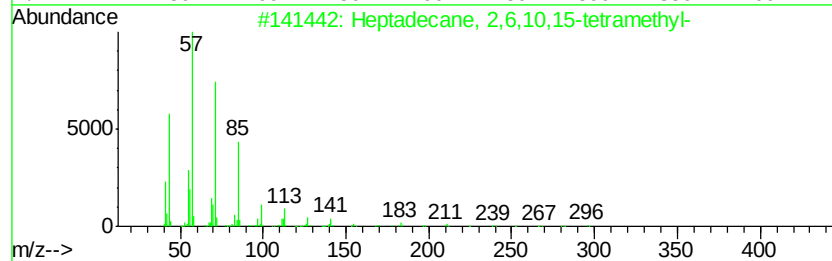
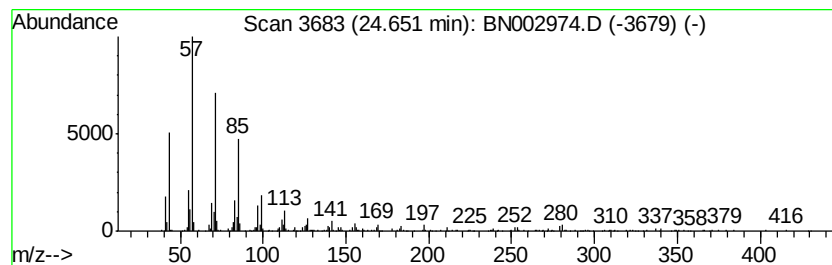
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.65	7.29 ng/ul	587871	Perylene-d12	23.44

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	93
2		Docosane	310	C22H46	000629-97-0	93
3		Triacontane	422	C30H62	000638-68-6	90
4		Hexacosane	366	C26H54	000630-01-3	90
5		Tridecane, 7-hexyl-	268	C19H40	007225-66-3	90



Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN100618\
Data File : BN002974.D
Acq On : 05 Oct 2018 21:55
Operator : MJ/SJ
Sample : J5230-04
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
A3YY0

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-BN091418MA.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
2,4,7,9-Tetrameth...	13.16	3.9	ng/ul	333264	3	14.26	1691040	20.0
n-Hexadecanoic acid	17.93	2.1	ng/ul	217835	4	17.02	2079090	20.0
unknown-01	18.09	2.4	ng/ul	245442	4	17.02	2079090	20.0
Phenanthrene, 2,5...	18.82	2.6	ng/ul	272967	4	17.02	2079090	20.0
Cyclopenta(def)ph...	18.97	2.5	ng/ul	256551	4	17.02	2079090	20.0
Pyrene, 1-methyl-	19.78	2.0	ng/ul	427465	5	21.23	4209780	20.0
11H-Benzo[a]fluor...	20.71	2.5	ng/ul	519301	5	21.23	4209780	20.0
unknown-02	20.94	4.4	ng/ul	928291	5	21.23	4209780	20.0
(DEL) Alkane: Str...	22.26	2.4	ng/ul	501668	5	21.23	4209780	20.0
Benzo[e]pyrene	22.94	4.2	ng/ul	341175	6	23.44	1613070	20.0
(DEL) Alkane: Str...	23.93	7.7	ng/ul	618921	6	23.44	1613070	20.0
(DEL) Alkane: Str...	24.65	7.3	ng/ul	587871	6	23.44	1613070	20.0