

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
 Data File : BN014864.D
 Acq On : 10 Jun 2021 14:36
 Operator : CG/JU
 Sample : M2618-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BG5R6

Integration Parameters: LSCINT.P

Integrator: RTE

Smoother: 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\SFAM-EPA-BN052521.M
 Title : SVOA CALIBRATION

Signal : TIC: BN014864.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.217	18	22	29	rBB4	25831	46591	1.02%	0.064%
2	3.840	122	128	132	rBB2	44104	60239	1.32%	0.083%
3	4.828	291	296	301	rBB	68278	88491	1.94%	0.121%
4	6.869	633	643	651	rBV	684724	1090473	23.85%	1.497%
5	7.034	663	671	679	rBB	444747	706067	15.44%	0.969%
6	7.228	698	704	713	rVB	712287	1131010	24.73%	1.552%
7	7.693	777	783	790	rBV	794654	1284831	28.10%	1.763%
8	8.399	895	903	917	rBV2	879366	1466516	32.07%	2.013%
9	8.846	973	979	986	rBB	677631	1078336	23.58%	1.480%
10	9.569	1093	1102	1111	rBV	544269	934189	20.43%	1.282%
11	10.105	1186	1193	1202	rBV2	943841	1599333	34.97%	2.195%
12	10.481	1247	1257	1264	rBV	1015172	1756817	38.42%	2.411%
13	10.616	1273	1280	1290	rBB2	713639	1266838	27.70%	1.739%
14	13.246	1722	1727	1734	rBB5	35952	74566	1.63%	0.102%
15	13.757	1808	1814	1820	rBB	1644452	2253804	49.29%	3.093%
16	14.040	1851	1862	1871	rBV	1770871	2758724	60.33%	3.786%
17	14.346	1907	1914	1923	rBV	1651192	2484750	54.34%	3.410%
18	14.557	1944	1950	1961	rBB	403686	697258	15.25%	0.957%
19	15.345	2077	2084	2090	rBV	2405764	3430913	75.03%	4.709%
20	15.463	2097	2104	2111	rBB5	138048	250633	5.48%	0.344%
21	15.587	2118	2125	2128	rBB8	30090	49922	1.09%	0.069%
22	17.098	2376	2382	2387	rBB	2024045	2827622	61.83%	3.881%
23	17.139	2387	2389	2393	rBB	150395	158749	3.47%	0.218%
24	17.198	2393	2399	2410	rBB2	2563153	3714723	81.23%	5.098%
25	18.016	2533	2538	2545	rBB9	74311	152119	3.33%	0.209%
26	18.486	2614	2618	2620	rBB5	31783	47717	1.04%	0.065%
27	19.169	2729	2734	2737	rBB	386366	473313	10.35%	0.650%
28	19.304	2753	2757	2758	rBB4	38915	43401	0.95%	0.060%
29	19.504	2786	2791	2800	rBB2	2914190	4572868	100.00%	6.276%
30	19.616	2807	2810	2812	rBB4	32563	44911	0.98%	0.062%
31	19.704	2821	2825	2828	rBB6	67137	68819	1.50%	0.094%
32	19.745	2828	2832	2834	rBB5	38399	49396	1.08%	0.068%
33	19.851	2847	2850	2853	rBB5	38563	56176	1.23%	0.077%
34	19.933	2863	2864	2867	rBB3	53628	44810	0.98%	0.062%
35	20.033	2869	2881	2883	rBB3	118391	182591	3.99%	0.251%
36	20.128	2895	2897	2899	rBB3	56153	47776	1.04%	0.066%

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37	20.186	2903	2907	2910	rBB6	72078	111265	2.43%	0.153%
38	20.292	2923	2925	2928	rBB4	71081	64492	1.41%	0.089%
39	20.416	2943	2946	2947	rBB3	97117	101318	2.22%	0.139%
40	20.445	2947	2951	2952	rBB4	116011	98896	2.16%	0.136%
41	20.586	2970	2975	2979	rBB8	72935	123975	2.71%	0.170%
42	20.663	2983	2988	2991	rBB7	138268	192579	4.21%	0.264%
43	20.704	2991	2995	2996	rBB4	43109	54227	1.19%	0.074%
44	20.780	3003	3008	3009	rBB5	97751	113639	2.49%	0.156%
45	20.798	3009	3011	3014	rBB4	46402	59124	1.29%	0.081%
46	20.875	3021	3024	3027	rBB5	52096	62461	1.37%	0.086%
47	20.992	3041	3044	3045	rBB3	55495	63625	1.39%	0.087%
48	21.039	3048	3052	3055	rBB6	123814	184779	4.04%	0.254%
49	21.104	3061	3063	3065	rBB3	113365	74500	1.63%	0.102%
50	21.122	3065	3066	3071	rBB5	127332	129120	2.82%	0.177%
51	21.228	3076	3084	3088	rBB5	270768	425986	9.32%	0.585%
52	21.304	3089	3097	3101	rBB	2121050	3073781	67.22%	4.219%
53	21.333	3101	3102	3105	rBB	94855	52356	1.14%	0.072%
54	21.375	3108	3109	3113	rBB4	44109	45750	1.00%	0.063%
55	21.410	3113	3115	3116	rBB2	61857	57979	1.27%	0.080%
56	21.457	3120	3123	3124	rBB3	74776	66975	1.46%	0.092%
57	21.539	3133	3137	3138	rBB4	56277	83387	1.82%	0.114%
58	21.551	3138	3139	3141	rBB2	139105	67700	1.48%	0.093%
59	21.633	3149	3153	3155	rBB5	86621	102214	2.24%	0.140%
60	21.722	3163	3168	3171	rBB7	187304	210705	4.61%	0.289%
61	21.769	3171	3176	3177	rBB5	49409	73490	1.61%	0.101%
62	21.804	3177	3182	3184	rBB6	167410	195553	4.28%	0.268%
63	21.839	3184	3188	3189	rBB4	66165	97635	2.14%	0.134%
64	21.886	3193	3196	3197	rBB3	97537	64190	1.40%	0.088%
65	21.922	3197	3202	3204	rBB6	83879	125427	2.74%	0.172%
66	21.998	3213	3215	3217	rBB3	105899	102758	2.25%	0.141%
67	22.027	3217	3220	3222	rBB4	130472	135171	2.96%	0.186%
68	22.080	3226	3229	3231	rBB4	62335	51261	1.12%	0.070%
69	22.151	3237	3241	3244	rBB5	188000	217692	4.76%	0.299%
70	22.239	3255	3256	3258	rBB2	92345	69067	1.51%	0.095%
71	22.280	3260	3263	3267	rBB6	94498	118232	2.59%	0.162%
72	22.375	3276	3279	3280	rBB3	108968	115014	2.52%	0.158%
73	22.439	3287	3290	3292	rBB4	68964	73211	1.60%	0.100%
74	22.486	3292	3298	3301	rBB6	166022	348885	7.63%	0.479%
75	22.516	3301	3303	3308	rBB6	112847	105461	2.31%	0.145%
76	22.557	3308	3310	3314	rBB5	85342	87485	1.91%	0.120%

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77	22.610	3317	3319	3320	rBB2	90206	50467	1.10%	0.069%
78	22.633	3322	3323	3324	rBB	49996	45794	1.00%	0.063%
79	22.663	3324	3328	3329	rBB4	111705	114894	2.51%	0.158%
80	22.710	3334	3336	3338	rBB3	68967	47259	1.03%	0.065%
81	22.745	3340	3342	3346	rBB5	75704	85639	1.87%	0.118%
82	22.798	3346	3351	3356	rBB9	180715	403076	8.81%	0.553%
83	22.892	3356	3367	3373	rBB9	473347	1432132	31.32%	1.966%
84	22.945	3373	3376	3381	rBB7	246915	339689	7.43%	0.466%
85	22.980	3381	3382	3385	rBB3	68439	54740	1.20%	0.075%
86	23.016	3387	3388	3391	rBB3	73127	43457	0.95%	0.060%
87	23.080	3391	3399	3405	rBB3	359519	841357	18.40%	1.155%
88	23.139	3407	3409	3410	rBB2	53004	61525	1.35%	0.084%
89	23.186	3414	3417	3419	rBB4	106572	121964	2.67%	0.167%
90	23.210	3419	3421	3427	rBB7	127325	206681	4.52%	0.284%
91	23.292	3430	3435	3436	rBB5	41772	192452	4.21%	0.264%
92	23.369	3444	3448	3449	rBB4	129365	125080	2.74%	0.172%
93	23.422	3451	3457	3462	rBB	1579125	2724955	59.59%	3.740%
94	23.457	3462	3463	3473	rBB8	78336	327570	7.16%	0.450%
95	23.569	3474	3482	3489	rBB	1527724	3218259	70.38%	4.417%
96	23.627	3489	3492	3493	rBB3	82240	74081	1.62%	0.102%
97	23.657	3496	3497	3499	rBB2	114898	45957	1.00%	0.063%
98	23.745	3509	3512	3513	rBB3	61508	52167	1.14%	0.072%
99	23.780	3513	3518	3527	rBB10	433109	912579	19.96%	1.253%
100	23.880	3527	3535	3538	rBB9	172972	384920	8.42%	0.528%
101	23.910	3538	3540	3541	rBB2	47727	45962	1.01%	0.063%
102	23.927	3541	3543	3547	rBB3	270828	332816	7.28%	0.457%
103	24.027	3556	3560	3561	rBB4	138264	163693	3.58%	0.225%
104	24.057	3564	3565	3572	rBB7	107544	62216	1.36%	0.085%
105	24.116	3572	3575	3581	rBB8	263656	418320	9.15%	0.574%
106	24.192	3586	3588	3593	rBB6	77237	64773	1.42%	0.089%
107	24.239	3593	3596	3600	rBB6	123016	164204	3.59%	0.225%
108	24.339	3602	3613	3623	rBB8	614919	1938563	42.39%	2.661%
109	24.445	3630	3631	3636	rBB5	78624	88223	1.93%	0.121%
110	24.492	3638	3639	3640	rBB	60381	52961	1.16%	0.073%
111	24.539	3640	3647	3648	rBB7	67872	167747	3.67%	0.230%
112	24.580	3651	3654	3655	rBB3	42252	50863	1.11%	0.070%
113	24.645	3659	3665	3675	rBB3	789078	1872663	40.95%	2.570%
114	24.751	3681	3683	3685	rBB3	91944	48699	1.06%	0.067%
115	24.786	3687	3689	3691	rBB3	66100	44298	0.97%	0.061%
116	24.816	3691	3694	3696	rBB4	85720	107995	2.36%	0.148%

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Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN052521.M
 Title : SVOA CALIBRATION

117	24.833	3696	3697	3700	rBB3	102566	93931	2.05%	0.129%
118	24.886	3700	3706	3712	rBB9	239293	528969	11.57%	0.726%
119	24.963	3718	3719	3724	rBV5	108749	120982	2.65%	0.166%
120	25.021	3727	3729	3731	rBB3	81658	64382	1.41%	0.088%
121	25.051	3733	3734	3735	rBB	139311	55393	1.21%	0.076%
122	25.098	3738	3742	3749	rBB8	240155	505084	11.05%	0.693%
123	25.204	3753	3760	3761	rBB7	41615	193020	4.22%	0.265%
124	25.327	3774	3781	3794	rBB4	634621	1720489	37.62%	2.361%
125	25.439	3798	3800	3802	rBB3	95876	83205	1.82%	0.114%
126	25.457	3802	3803	3804	rBB	71938	42933	0.94%	0.059%
127	25.480	3804	3807	3809	rBB4	50752	65730	1.44%	0.090%
128	25.580	3820	3824	3827	rBB6	89726	96095	2.10%	0.132%
129	25.704	3842	3845	3849	rBB6	79171	89544	1.96%	0.123%
130	25.768	3849	3856	3868	rBB6	322262	884435	19.34%	1.214%
131	25.880	3873	3875	3876	rBB2	109875	55413	1.21%	0.076%
132	25.916	3876	3881	3885	rBB8	161656	246293	5.39%	0.338%
133	26.033	3899	3901	3903	rBB3	87260	83404	1.82%	0.114%
134	26.068	3906	3907	3909	rBB2	73111	51199	1.12%	0.070%
135	26.104	3911	3913	3918	rBB6	126476	87229	1.91%	0.120%
136	26.145	3918	3920	3923	rBB4	65568	45678	1.00%	0.063%
137	26.210	3923	3931	3940	rBB6	592153	1652927	36.15%	2.269%
138	26.321	3945	3950	3951	rBB5	52771	160872	3.52%	0.221%
139	26.474	3973	3976	3977	rBB3	42106	58826	1.29%	0.081%
140	26.539	3985	3987	3988	rBB2	71786	75796	1.66%	0.104%
141	26.557	3988	3990	3994	rBB5	99906	124266	2.72%	0.171%
142	26.668	4005	4009	4010	rBB4	55742	79962	1.75%	0.110%
143	26.692	4010	4013	4016	rBB5	152294	173108	3.79%	0.238%
144	26.792	4027	4030	4031	rBB3	62712	86057	1.88%	0.118%
145	26.851	4039	4040	4044	rBB4	48822	43397	0.95%	0.060%
146	26.886	4044	4046	4049	rBV4	76392	73647	1.61%	0.101%
147	27.004	4061	4066	4071	rBB8	242178	513486	11.23%	0.705%
148	27.110	4083	4084	4086	rBB2	95154	62843	1.37%	0.086%
149	27.151	4089	4091	4092	rBB2	79044	54488	1.19%	0.075%
150	27.192	4095	4098	4102	rBB6	193091	152412	3.33%	0.209%
151	27.268	4108	4111	4113	rBB4	72119	57028	1.25%	0.078%
152	27.345	4123	4124	4128	rBB4	99052	57813	1.26%	0.079%
153	27.386	4128	4131	4132	rBB3	134906	85094	1.86%	0.117%
154	27.427	4136	4138	4141	rBB4	89386	107653	2.35%	0.148%
155	27.457	4141	4143	4146	rBB4	191996	138086	3.02%	0.190%
156	27.533	4155	4156	4158	rBB2	92913	57956	1.27%	0.080%

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157	27.639	4170	4174	4175	rBB4	81163	50392	1.10%	0.069%
158	27.662	4175	4178	4180	rBB4	55949	52607	1.15%	0.072%
159	27.692	4180	4183	4186	rBB5	110770	109962	2.40%	0.151%
160	27.715	4186	4187	4192	rBB5	81896	116945	2.56%	0.161%
161	27.845	4208	4209	4213	rBB4	50491	54351	1.19%	0.075%
162	27.968	4227	4230	4233	rBV5	88315	121127	2.65%	0.166%
163	28.021	4233	4239	4240	rBB6	15548	62423	1.37%	0.086%
164	28.057	4242	4245	4246	rBB3	58945	68457	1.50%	0.094%
165	28.074	4246	4248	4251	rBB4	90285	84495	1.85%	0.116%
166	28.345	4291	4294	4296	rBB4	54500	61316	1.34%	0.084%
167	28.468	4313	4315	4318	rBB4	93468	77673	1.70%	0.107%
168	28.539	4325	4327	4330	rBB4	63437	62185	1.36%	0.085%
169	28.568	4331	4332	4334	rBB2	81670	55406	1.21%	0.076%
170	28.604	4334	4338	4343	rBB8	113689	140304	3.07%	0.193%
171	28.651	4343	4346	4348	rBB4	72289	89703	1.96%	0.123%
172	28.698	4351	4354	4355	rBB3	91309	59472	1.30%	0.082%
173	28.821	4372	4375	4379	rBB6	86813	94828	2.07%	0.130%
174	28.904	4386	4389	4391	rBB4	60492	50895	1.11%	0.070%
175	28.939	4394	4395	4398	rBB3	75661	52767	1.15%	0.072%
176	28.986	4401	4403	4404	rBB2	46830	46297	1.01%	0.064%
177	29.074	4412	4418	4419	rBB6	68997	101062	2.21%	0.139%
178	29.186	4436	4437	4440	rBB3	100382	80894	1.77%	0.111%
179	29.262	4449	4450	4453	rBB3	79190	64322	1.41%	0.088%
180	29.327	4459	4461	4463	rBB3	88321	46075	1.01%	0.063%

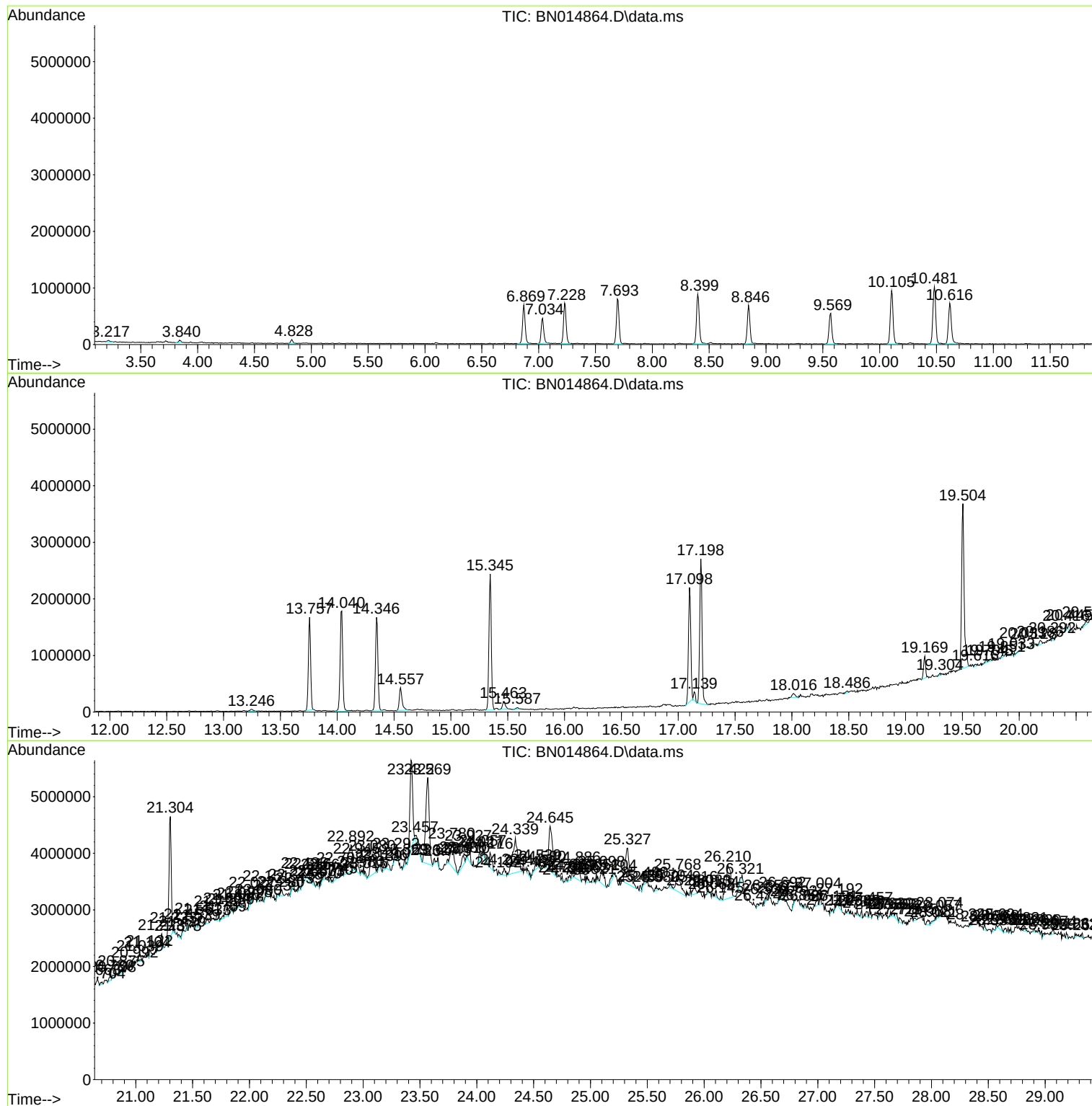
Sum of corrected areas: 72860430

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BG5R6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN052521.M
Quant Title : SVOA CALIBRATION

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



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Data File : BN014864.D
Acq On : 10 Jun 2021 14:36
Operator : CG/JU
Sample : M2618-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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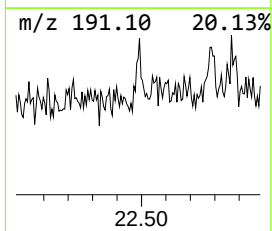
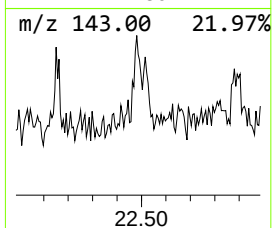
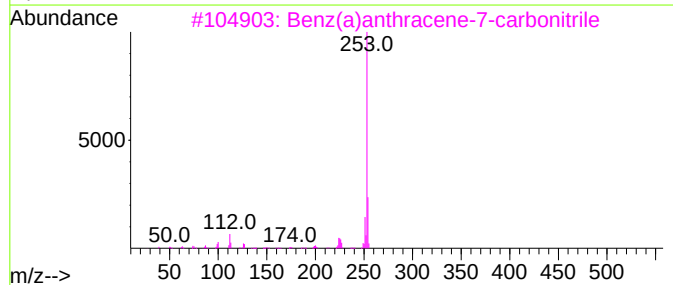
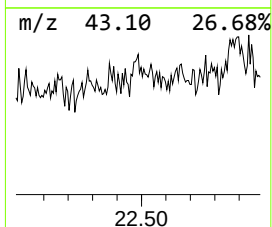
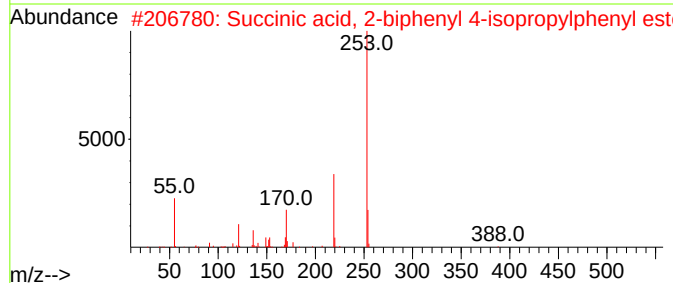
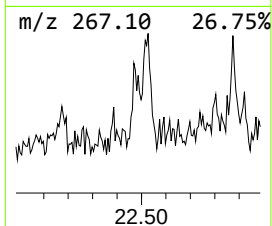
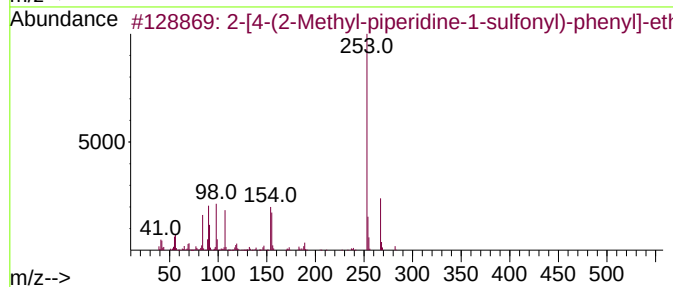
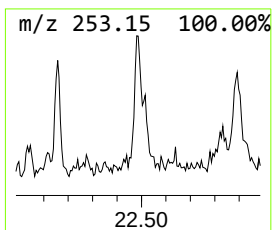
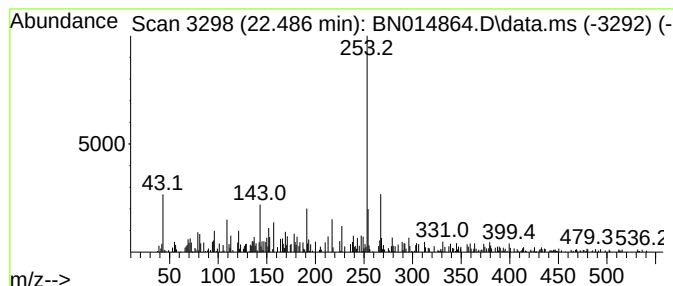
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Quant Title : SVOA CALIBRATION

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

Peak Number 1 2-[4-(2-Methyl-piperidine-1-... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.486	2.17 ng/ul	348885	Perylene-d12	23.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2	-	[4-(2-Methyl-piperidine-1-sulf...	282	C14H22N2O2S	1000295-08-5	59
2	Succinic acid, 2-biphenyl 4-isop...	388	C25H24O4	1000360-72-3	53		
3	Benz(a)anthracene-7-carbonitrile	253	C19H11N	007476-08-6	49		
4	Succinic acid, di(2-phenylphenyl...	422	C28H22O4	1000330-01-1	49		
5	1,3-Isoindolinedione, 2-(4-metho...	253	C15H11NO3	002142-04-3	47		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
Data File : BN014864.D
Acq On : 10 Jun 2021 14:36
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Sample : M2618-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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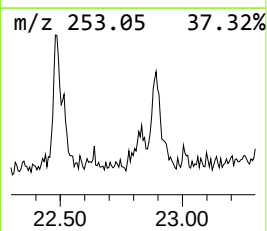
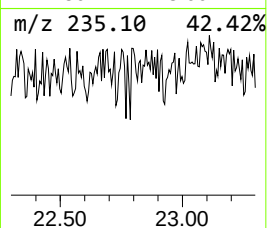
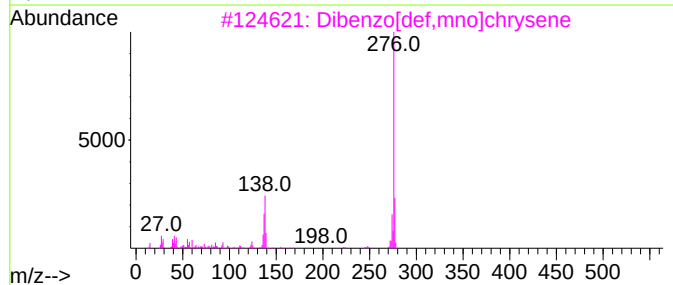
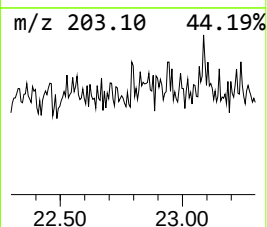
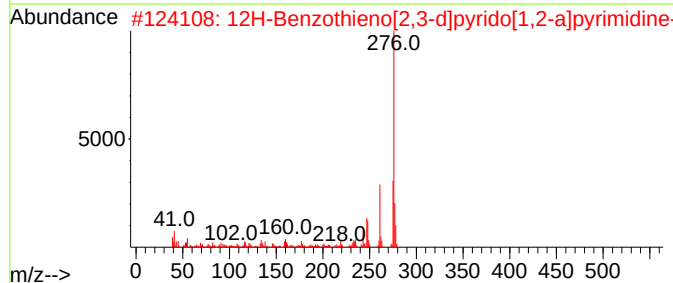
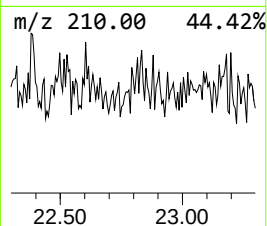
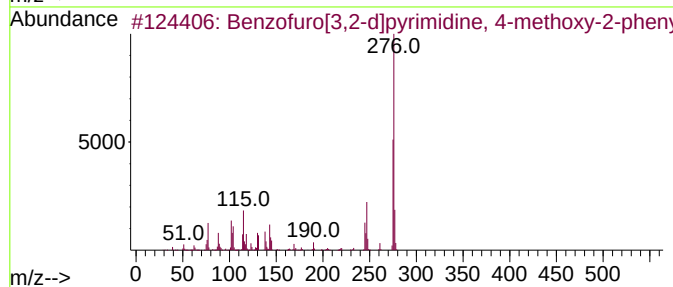
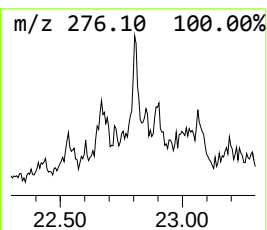
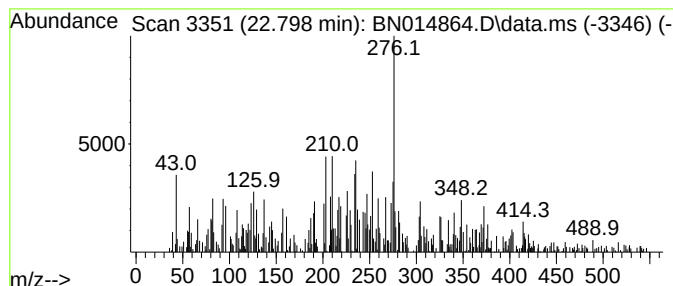
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Quant Title : SVOA CALIBRATION

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

Peak Number 2 unknown-01 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.798	2.50 ng/ul	403076	Perylene-d12	23.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzofuro[3,2-d]pyrimidine, 4-me...	276	C17H12N2O2	312265-36-4	30
2			12H-Benzothieno[2,3-d]pyrido[1,2...	276	C14H16N2S2	102254-63-7	27
3			Dibenzo[def,mno]chrysene	276	C22H12	000191-26-4	25
4			[1,3]Benzimidazo[2',1':2,3]thiaz...	276	C15H8N4S	1000319-10-6	25
5			1H-Pyrrol-5-amine, 4-(2-benzoxaz...	276	C16H12N4O	039116-39-7	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
Data File : BN014864.D
Acq On : 10 Jun 2021 14:36
Operator : CG/JU
Sample : M2618-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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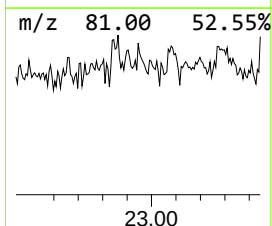
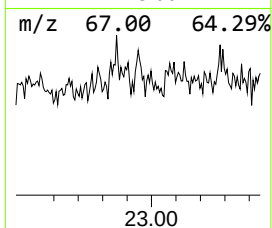
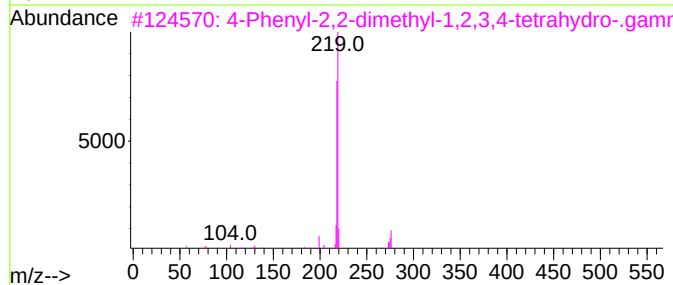
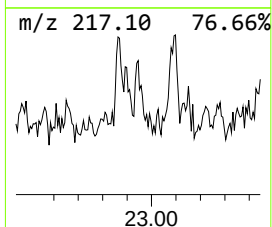
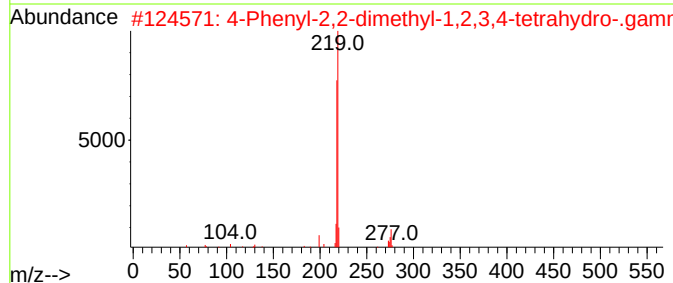
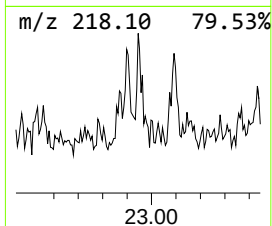
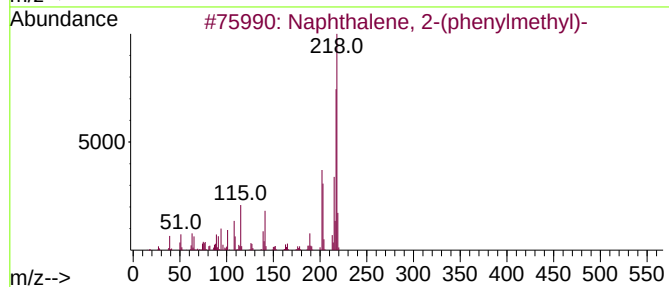
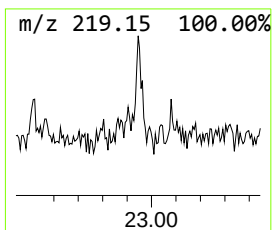
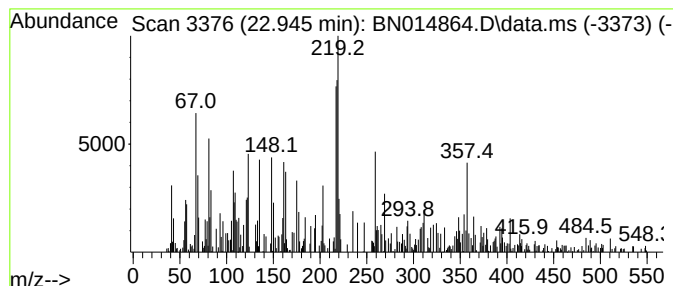
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-02 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.945	2.11 ng/ul	339689	Perylene-d12	23.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-(phenylmethyl)-	218	C17H14	000613-59-2	43
2			4-Phenyl-2,2-dimethyl-1,2,3,4-te...	276	C19H20N2	022315-73-7	38
3			4-Phenyl-2,2-dimethyl-1,2,3,4-te...	276	C19H20N2	022315-73-7	38
4			Quinoline, 6-methyl-2-phenyl-	219	C16H13N	027356-46-3	30
5			Quinoline, 2-(4-methylphenyl)-	219	C16H13N	024667-94-5	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
Data File : BN014864.D
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Operator : CG/JU
Sample : M2618-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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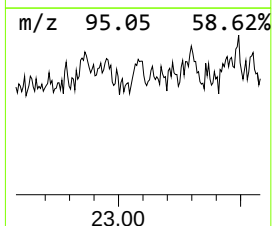
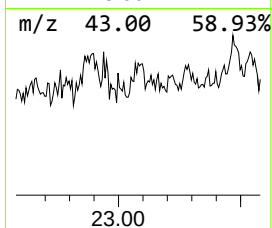
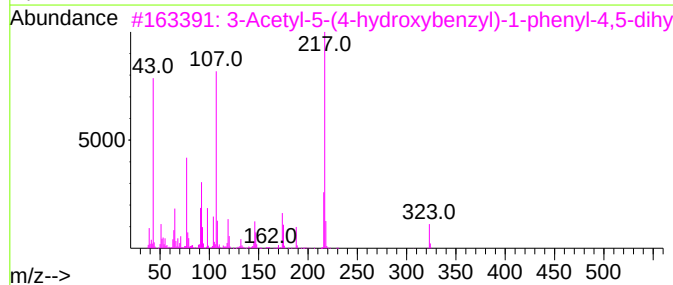
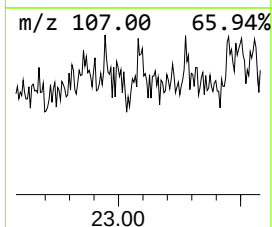
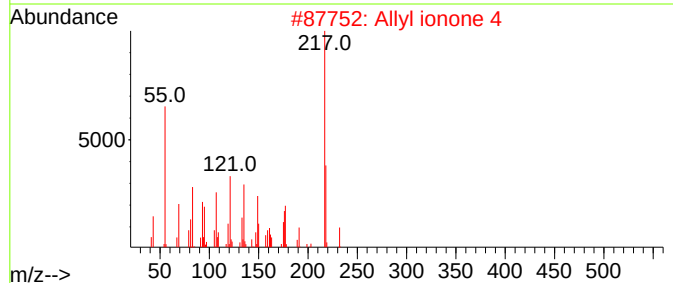
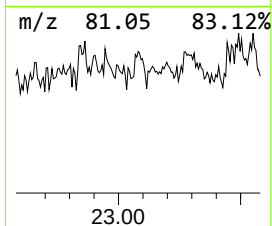
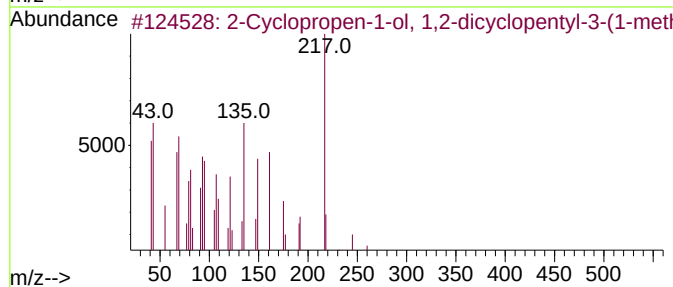
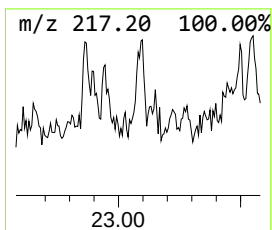
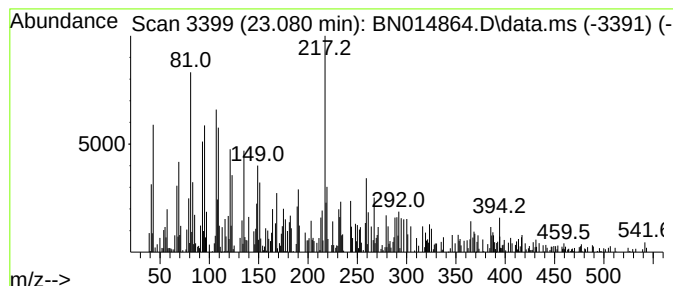
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-03 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.080	5.23 ng/ul	841357	Perylene-d12	23.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Cyclopropen-1-ol, 1,2-dicycloprop...	276	C18H28O2	069611-52-5	43		
2	Allyl ionone 4	232	C16H24O	1000284-93-0	38		
3	3-Acetyl-5-(4-hydroxybenzyl)-1-p...	323	C18H17N3O3	1000210-18-2	22		
4	3-Amino-6-ethyl-5-methyl-thieno[...	217	C11H11N3S	1000300-52-6	15		
5	7H-Benzo[c]carbazole	217	C16H11N	000205-25-4	14		



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
Data File : BN014864.D
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Sample : M2618-03
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ALS Vial : 6 Sample Multiplier: 1

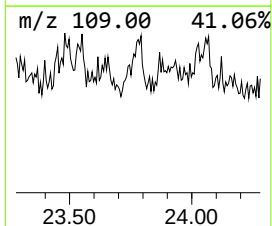
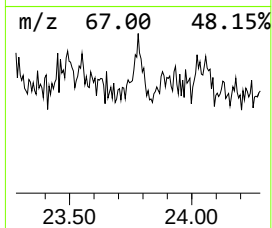
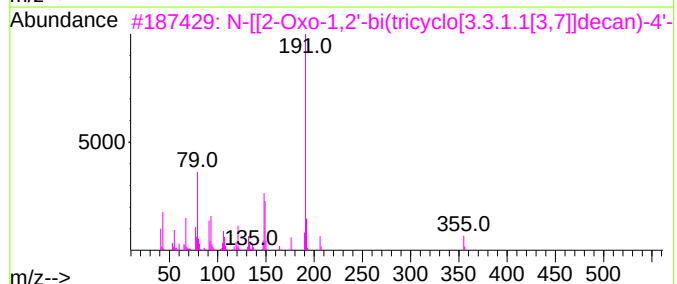
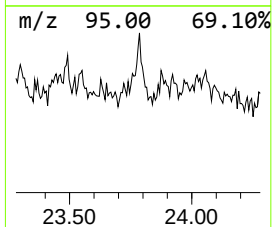
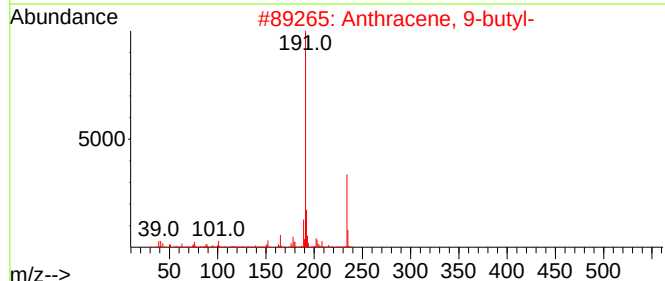
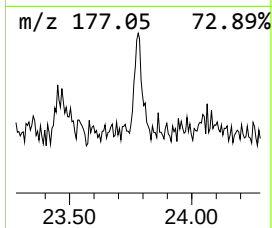
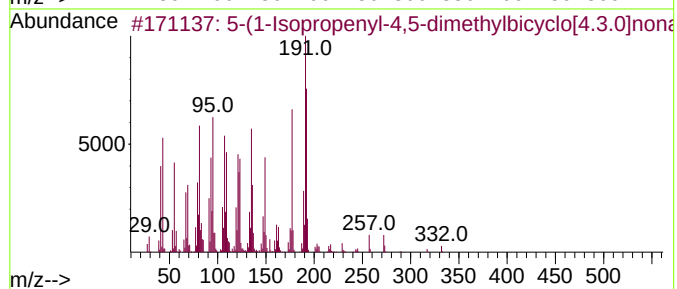
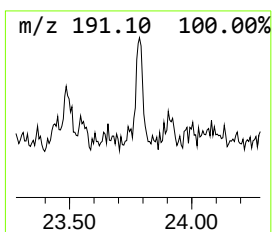
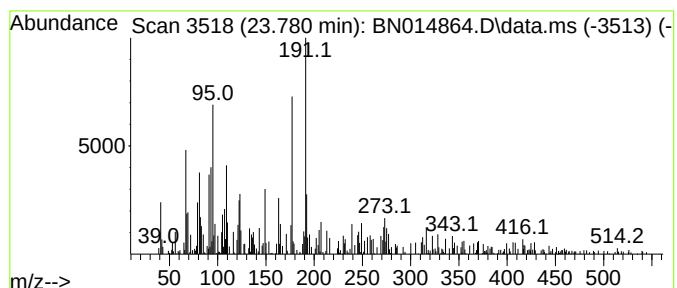
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 5-(1-Isopropenyl-4,5-dimeth... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.780	5.67 ng/ul	912579	Perylene-d12	23.569	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-(1-Isopropenyl-4,5-dimethylbic...	332	C22H36O2	1000195-69-8	52
2	Anthracene, 9-butyl-	234	C18H18	001498-69-7	25
3	N-[[2-Oxo-1,2'-bi(tricyclo[3.3.1...	355	C23H33NO2	1000305-39-6	25
4	1,2,3,4-Tetrahydropyrimidin-2,4-...	416	C21H28N4O5	1000127-97-3	25
5	Terephthalic acid, monoamide, N-...	311	C19H21NO3	1000323-93-8	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
Data File : BN014864.D
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Misc :
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Instrument :
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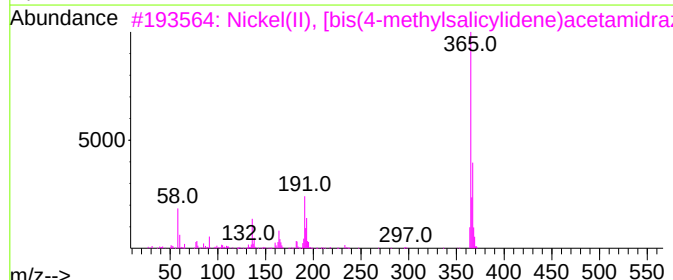
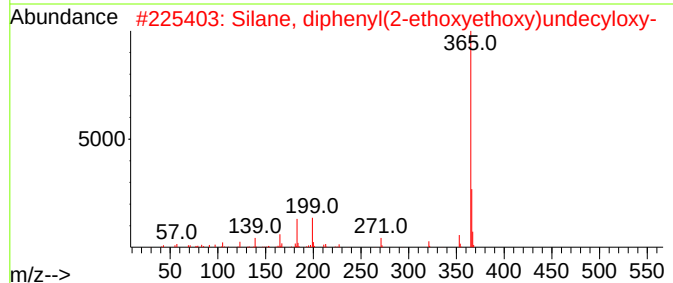
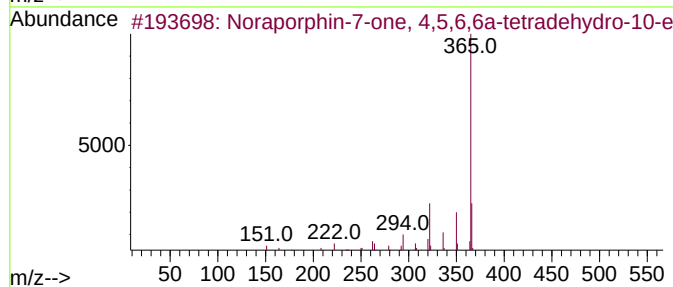
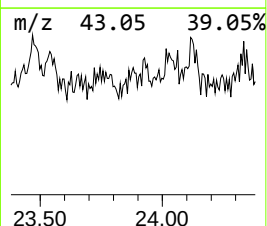
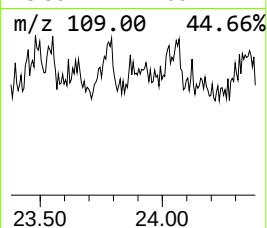
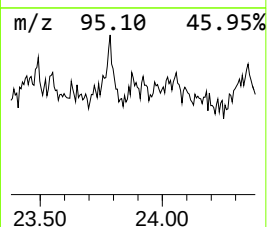
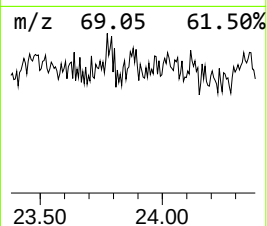
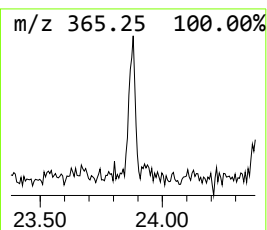
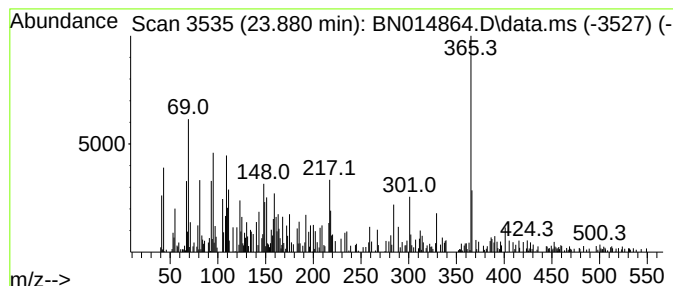
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-04 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.880	2.39 ng/ul	384920	Perylene-d12	23.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Noraporphin-7-one, 4,5,6,6a-tetr...	365	C21H19NO5	015358-01-7	38
2			Silane, diphenyl(2-ethoxyethoxy)...	442	C27H42O3Si	1000367-66-8	27
3			Nickel(II), [bis(4-methylsalicyl...	365	C18H17N3NiO2	1000164-61-6	27
4			Cobalt,bis[(1,2,3,4,5,6-.eta.)-1...	365	C22H20B2Co	036534-31-3	27
5			cis-11,14-Eicosadienoic acid, te...	422	C26H50O2Si	1000333-61-9	17



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
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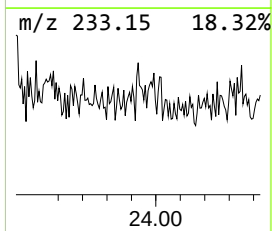
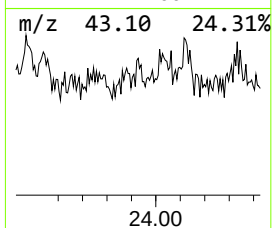
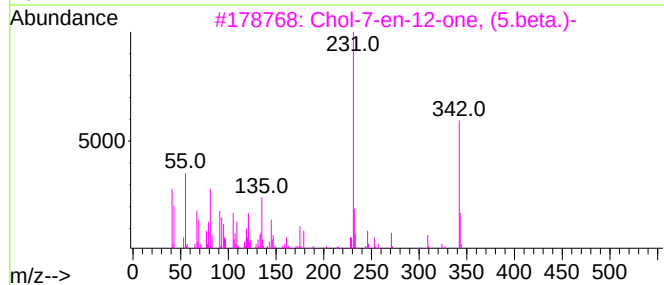
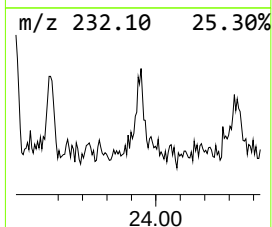
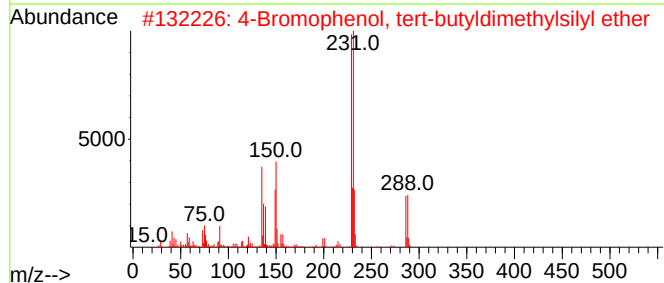
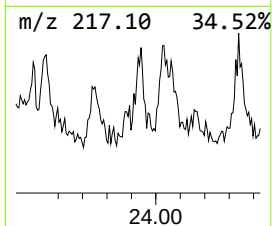
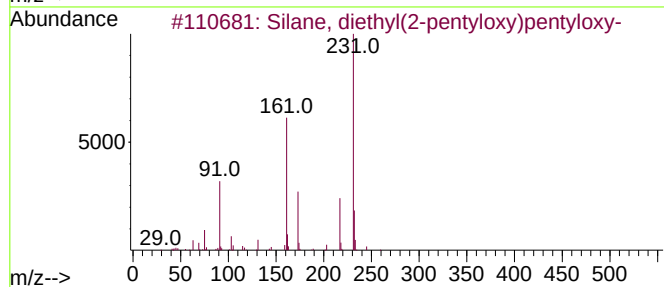
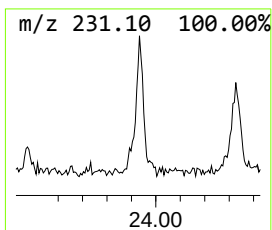
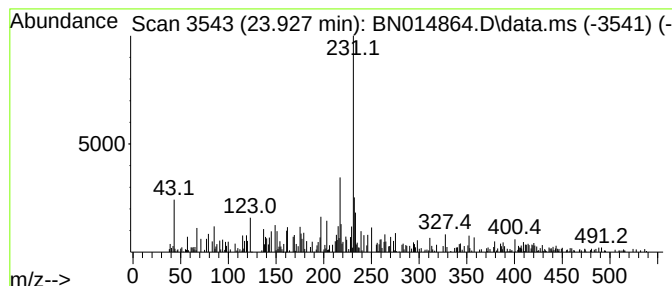
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Silane, diethyl(2-pentyloxy... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.927	2.07 ng/ul	332816	Perylene-d12	23.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Silane, diethyl(2-pentyloxy)pent...	260	C14H32O2Si	1000362-98-0	59
2			4-Bromophenol, tert-butyldimethy...	286	C12H19BrOSi	1000333-45-6	59
3			Chol-7-en-12-one, (5.beta.)-	342	C24H38O	054411-74-4	53
4			Pyrimidine, 2-dimethylamino-4,6-...	231	C12H13N3O2	1000261-69-3	53
5			Pyrimido[5,4-b]pyrido[3,2-d]thie...	231	C11H9N3OS	201681-22-3	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
Data File : BN014864.D
Acq On : 10 Jun 2021 14:36
Operator : CG/JU
Sample : M2618-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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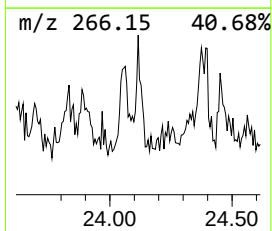
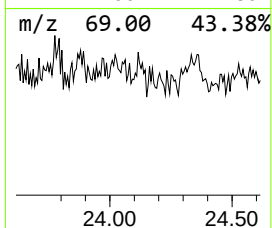
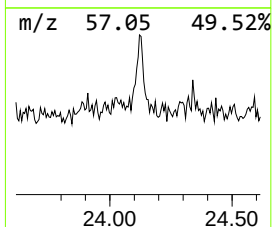
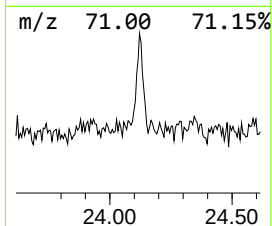
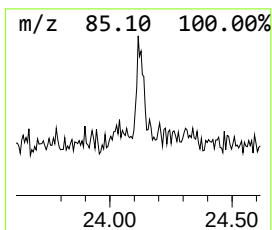
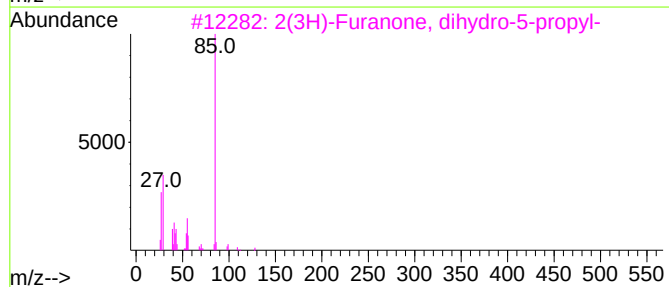
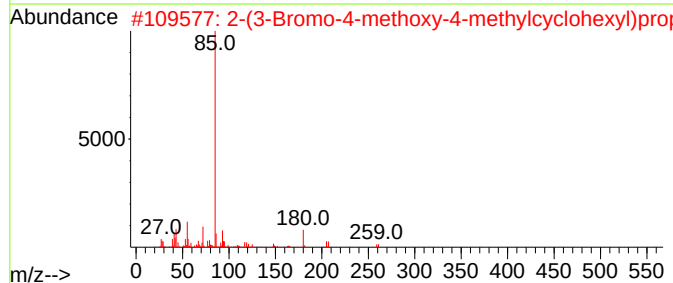
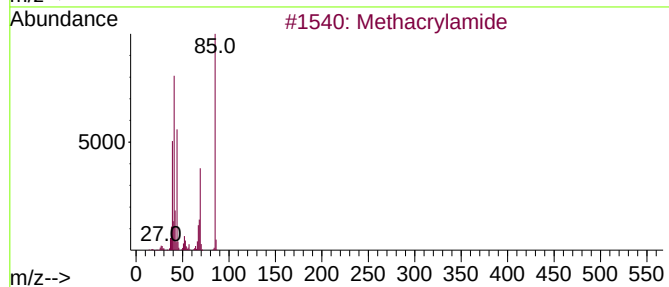
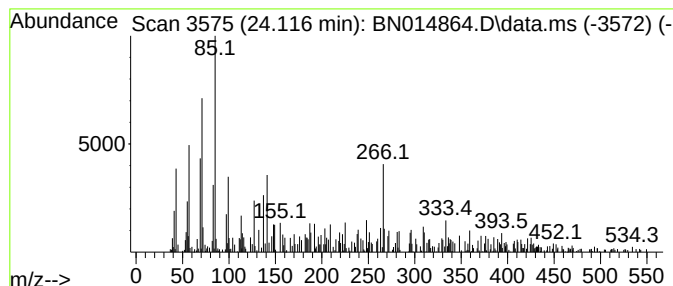
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-05 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.116	2.60 ng/ul	418320	Perylene-d12	23.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methacrylamide	85	C4H7NO	000079-39-0	46
2			2-(3-Bromo-4-methoxy-4-methylcyc...	259	C11H18BrNO	1000210-71-3	46
3			2(3H)-Furanone, dihydro-5-propyl-	128	C7H12O2	000105-21-5	43
4			Sulfurous acid, hexyl nonyl ester	292	C15H32O3S	1000309-13-1	43
5			Valeric acid, 4-nitrophenyl ester	223	C11H13NO4	1000307-98-9	43



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Sample : M2618-03
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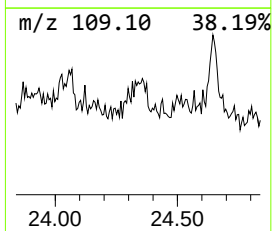
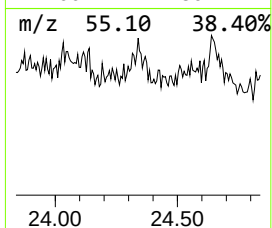
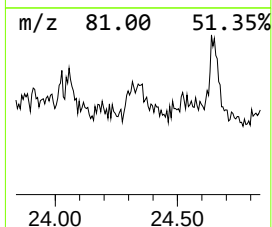
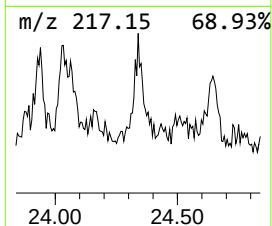
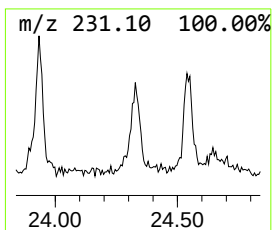
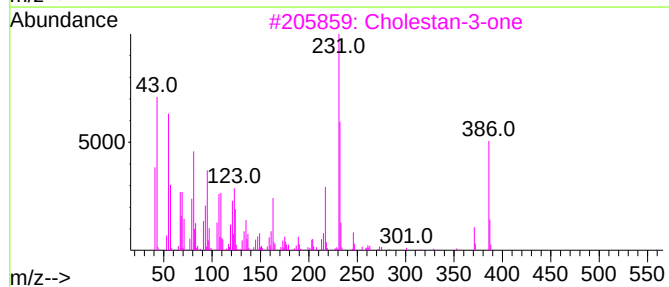
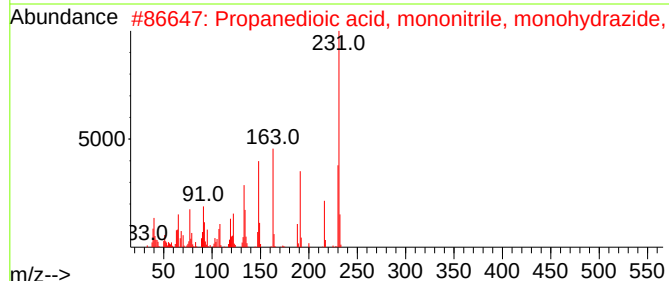
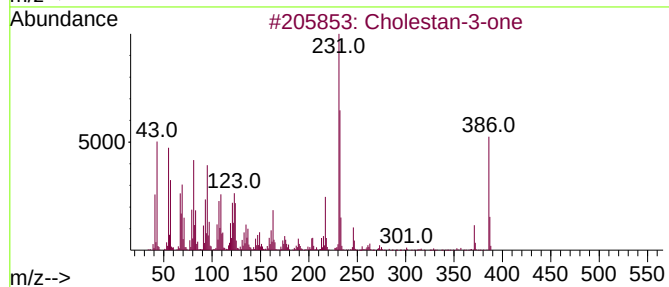
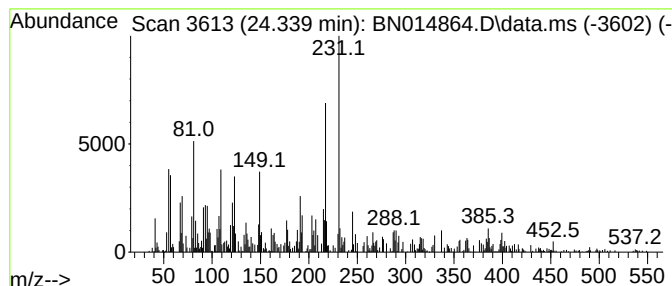
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Cholestan-3-one Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.339	12.05 ng/ul	1938560	Perylene-d12	23.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cholestan-3-one	386	C27H46O	015600-08-5	53
2			Propanedioic acid, mononitrile, ...	231	C12H13N3O2	004980-37-4	50
3			Cholestan-3-one	386	C27H46O	015600-08-5	47
4			Butylphosphonic acid, 2-bromo-4-...	422	C18H29BrF03P	1000315-12-4	30
5			2-Iodobenzoyl chloride	266	C7H4ClIO	000609-67-6	30



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Sample : M2618-03
Misc :
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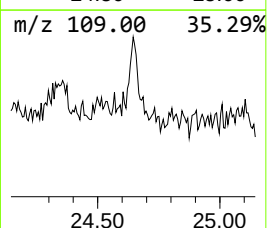
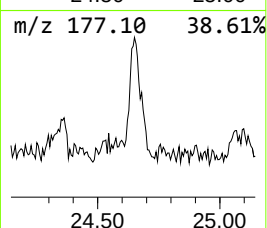
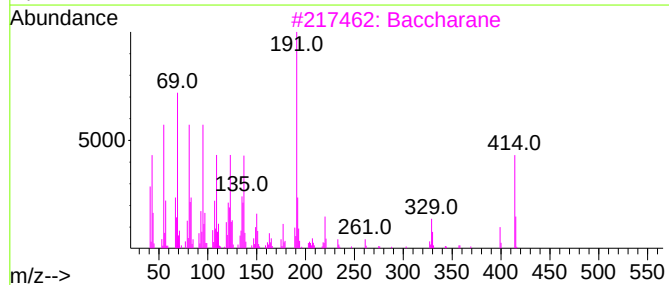
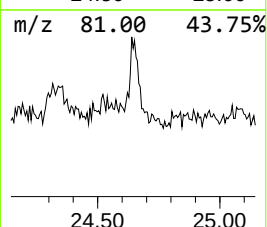
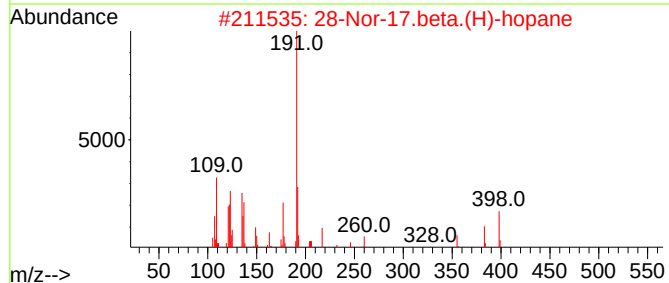
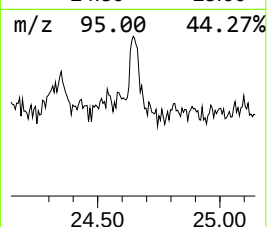
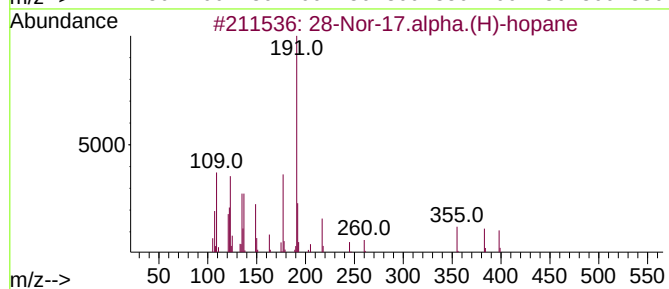
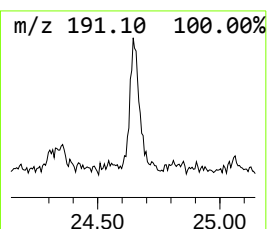
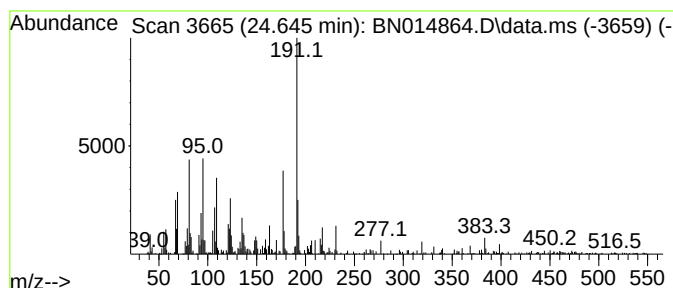
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 28-Nor-17.alpha.(H)-hopane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.645	11.64 ng/ul	1872660	Perylene-d12	23.569	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	28-Nor-17.alpha.(H)-hopane	398	C29H50	053584-60-4	86
2	28-Nor-17.beta.(H)-hopane	398	C29H50	036728-72-0	38
3	Baccharane	414	C30H54	036441-74-4	38
4	2,4,7-Pteridinetriamine, 6-methyl-	191	C7H9N7	017539-50-3	30
5	3-Buten-2-one, 4-(2,5,6,6-tetram...	206	C14H22O	000079-70-9	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
Data File : BN014864.D
Acq On : 10 Jun 2021 14:36
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Misc :
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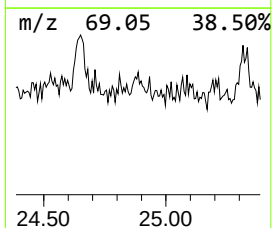
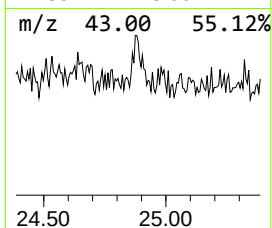
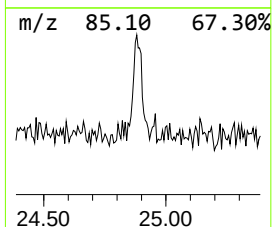
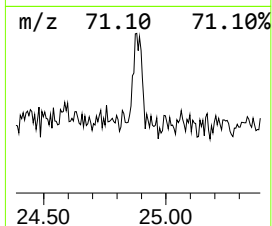
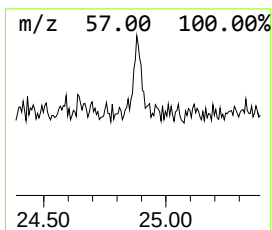
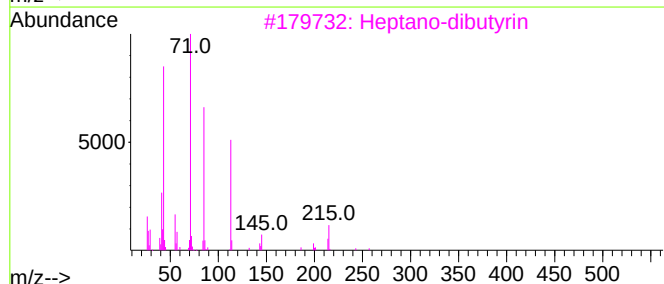
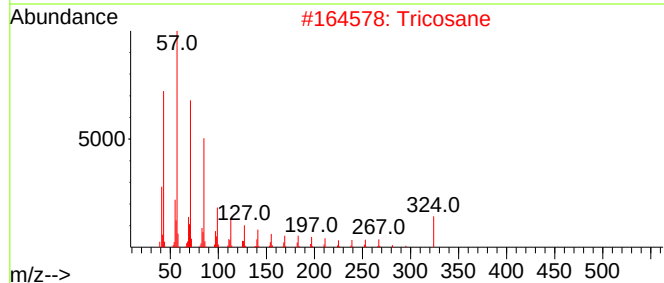
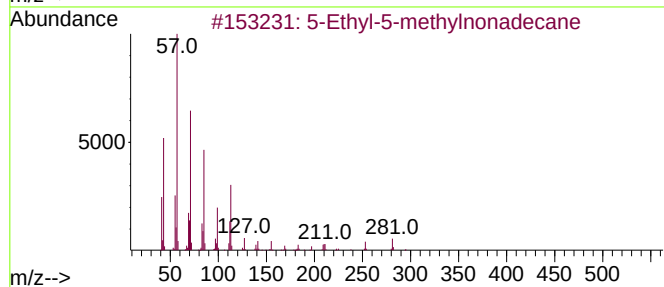
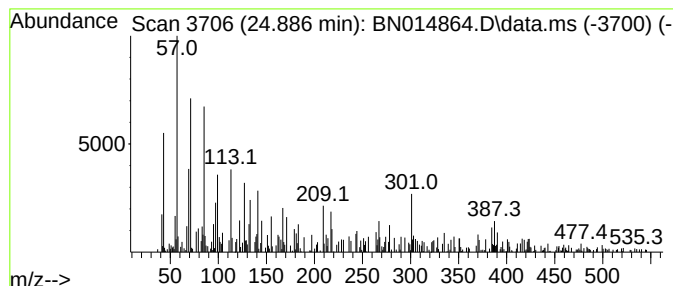
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.886	3.29 ng/ul	528969	Perylene-d12	23.569

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		5-Ethyl-5-methylnonadecane	310	C22H46	1000360-42-7	50
2		Tricosane	324	C23H48	000638-67-5	50
3		Heptano-dibutyryn	344	C18H32O6	065235-13-4	40
4		5-Undecanone	170	C11H22O	033083-83-9	40
5		Squalane	422	C30H62	000111-01-3	40



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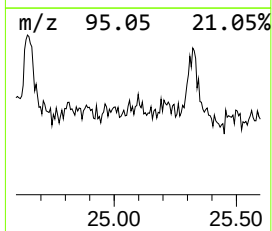
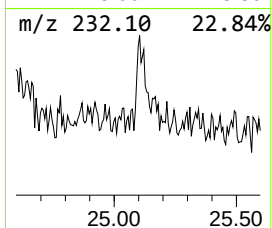
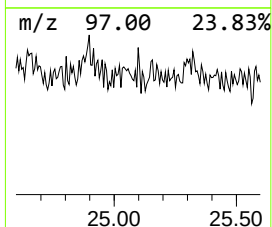
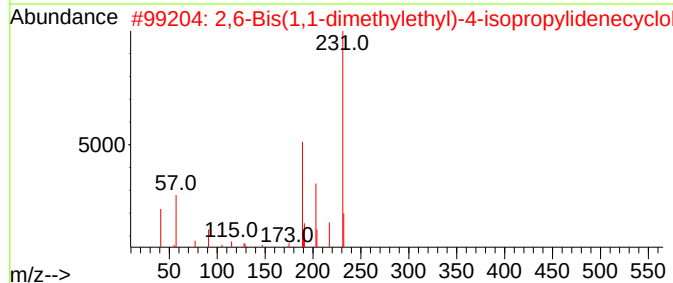
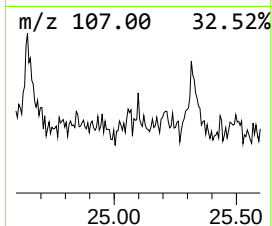
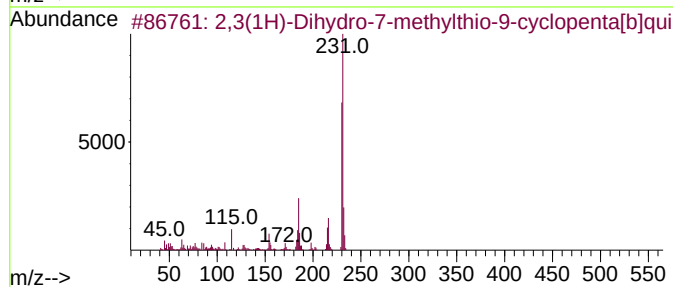
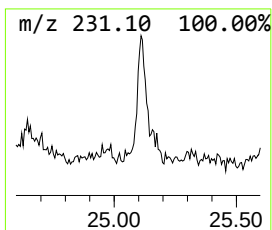
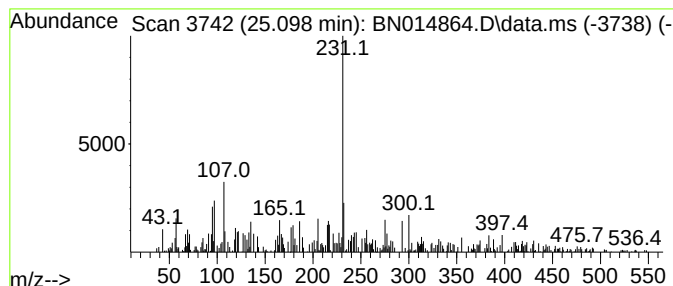
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 2,3(1H)-Dihydro-7-methylthi... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
25.098	3.14 ng/ul	505084	Perylene-d12	23.569	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3(1H)-Dihydro-7-methylthio-9-c...	231	C13H13NOS	1000257-08-5	72
2	2,6-Bis(1,1-dimethylethyl)-4-iso...	246	C17H26O	005427-02-1	64
3	Silane, diethylethoxy(2-ethylhex...	260	C14H32O2Si	1000363-04-5	64
4	10-Methyl-1,3,4,10-tetrahydro-1,...	231	C12H13N3S	026228-45-5	64
5	2-Cyano-3-phenylquinoxaline	231	C15H9N3	059393-45-2	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
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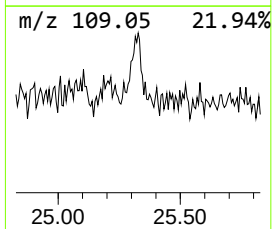
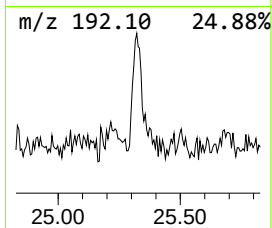
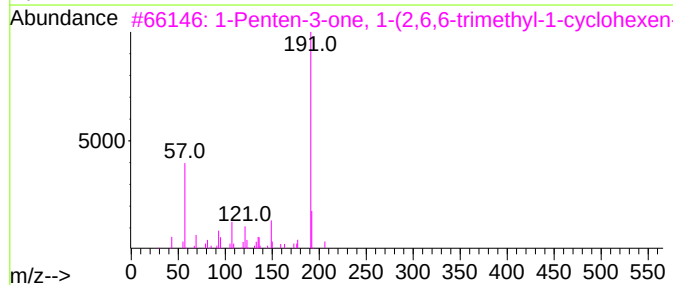
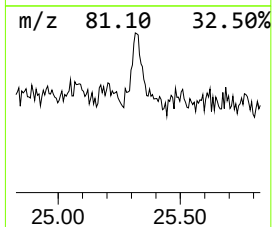
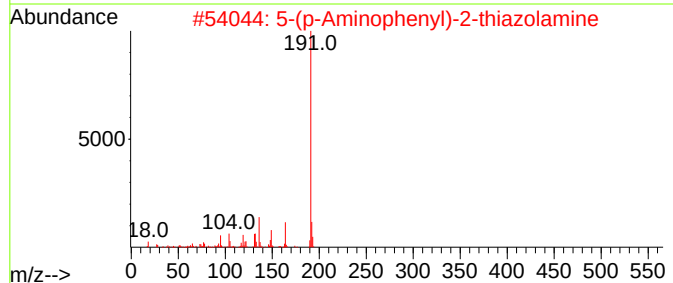
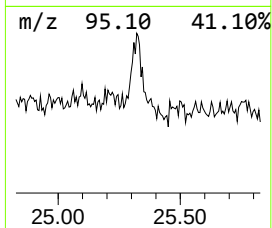
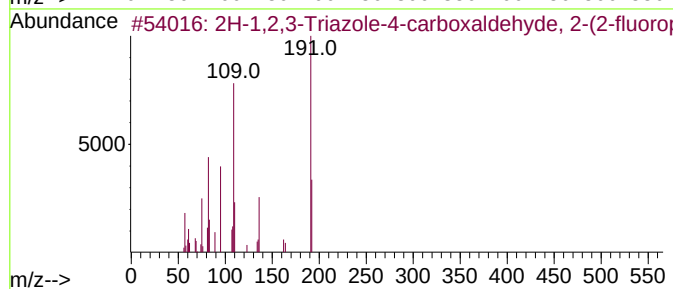
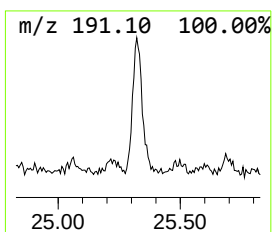
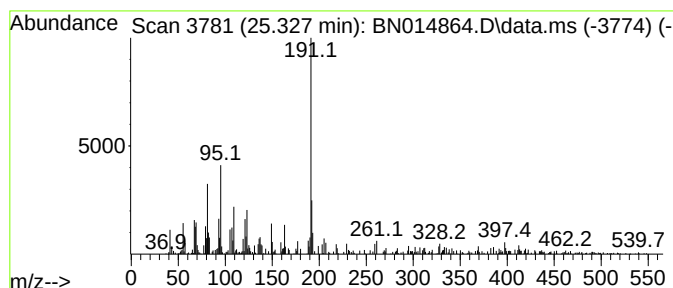
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 2H-1,2,3-Triazole-4-carboxa... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
25.327	10.69 ng/ul	1720490	Perylene-d12	23.569	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-1,2,3-Triazole-4-carboxaldehy...	191	C9H6FN3O	051306-43-5	53
2	5-(p-Aminophenyl)-2-thiazolamine	191	C9H9N3S	090349-87-4	38
3	1-Penten-3-one, 1-(2,6,6-trimeth...	206	C14H22O	000127-43-5	38
4	9H-Fluorene-9-carbonitrile	191	C14H9N	001529-40-4	35
5	9H-Fluorene-9-carbonitrile	191	C14H9N	001529-40-4	35



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Acq On : 10 Jun 2021 14:36
Operator : CG/JU
Sample : M2618-03
Misc :
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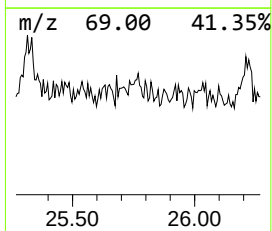
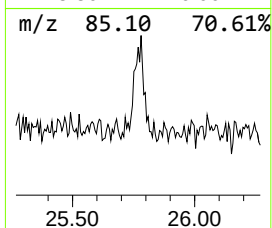
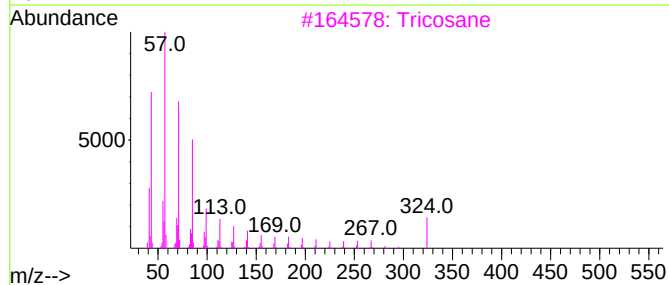
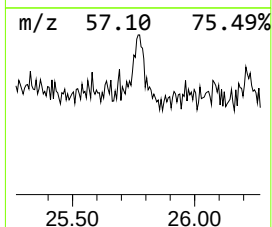
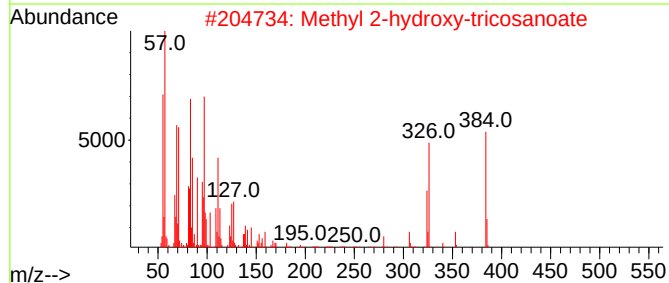
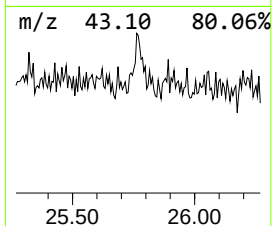
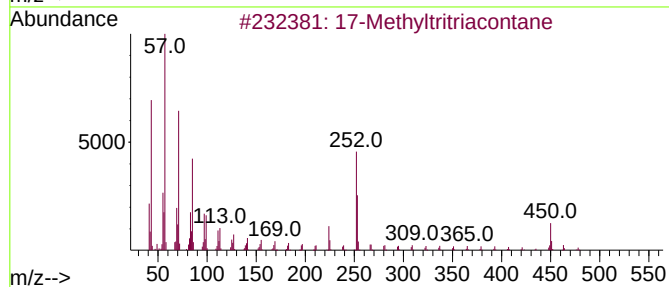
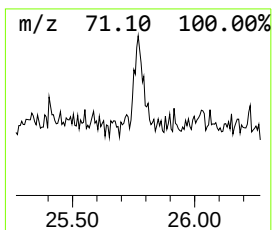
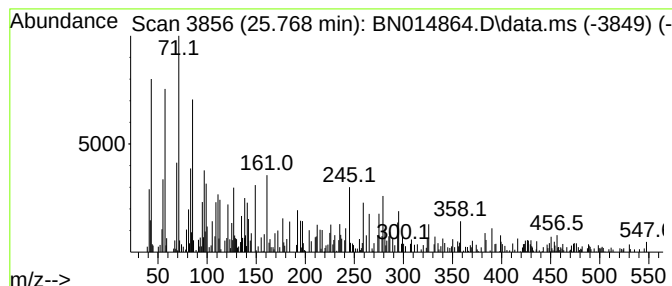
Instrument :
BNA_N
ClientSampleId :
BG5R6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN052521.M
Quant Title : SVOA CALIBRATION

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

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
25.769	5.50 ng/ul	884435	Perylene-d12	23.569	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	17-Methyltritriacontane	479	C34H70	058349-88-5	47
2	Methyl 2-hydroxy-tricosanoate	384	C24H48O3	1000336-21-4	45
3	Tricosane	324	C23H48	000638-67-5	38
4	3,3,13,13-Tetraethylpentadecane	324	C23H48	1000360-42-3	38
5	Octadecane, 3-ethyl-5-(2-ethylbu...	366	C26H54	055282-12-7	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
Data File : BN014864.D
Acq On : 10 Jun 2021 14:36
Operator : CG/JU
Sample : M2618-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

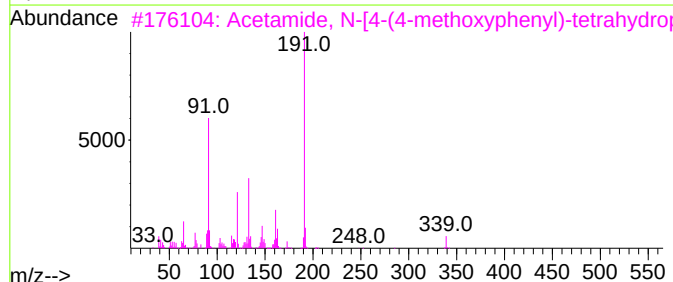
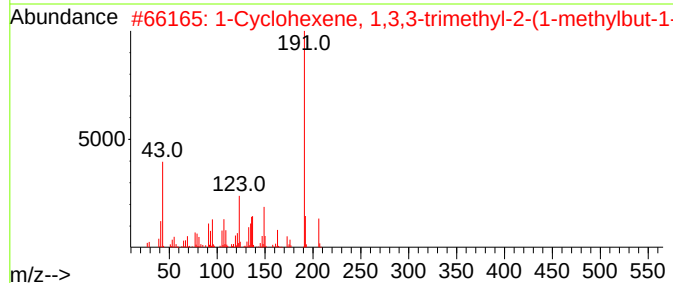
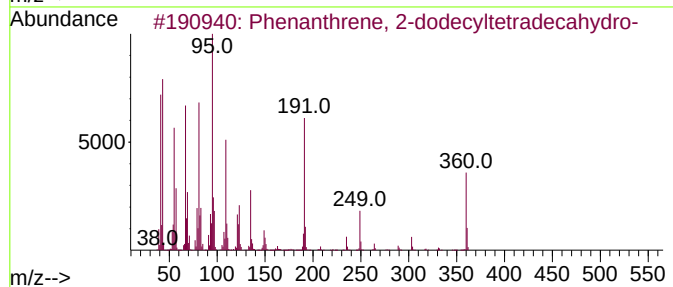
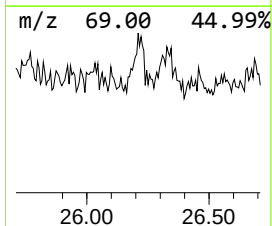
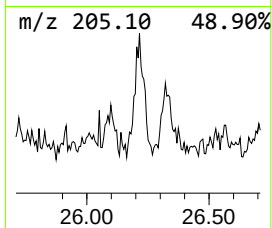
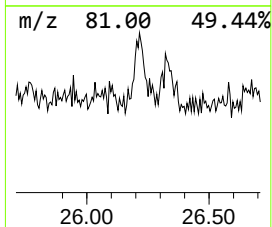
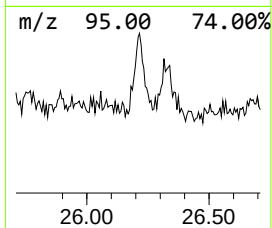
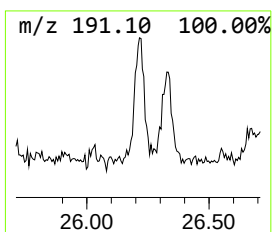
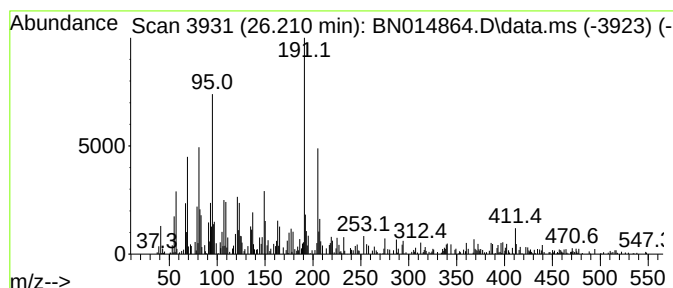
Instrument :
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN052521.M
Quant Title : SVOA CALIBRATION

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

Peak Number 15 Phenanthrene, 2-dodecyltetra... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
26.210	10.27 ng/ul	1652930	Perylene-d12	23.569	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2-dodecyltetradeca...	360	C26H48	055334-22-0	62
2	1-Cyclohexene, 1,3,3-trimethyl-2...	206	C14H22O	1000197-08-4	38
3	Acetamide, N-[4-(4-methoxyphenyl)...]	339	C21H25NO3	1000305-72-2	30
4	2(3H)-Thiazolone, 4-(4-methylphe...	191	C10H9NOS	1000338-15-1	25
5	2,4,5,5,8a-Pentamethyl-6,7,8,8a-...	206	C14H22O	1000195-40-9	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
Data File : BN014864.D
Acq On : 10 Jun 2021 14:36
Operator : CG/JU
Sample : M2618-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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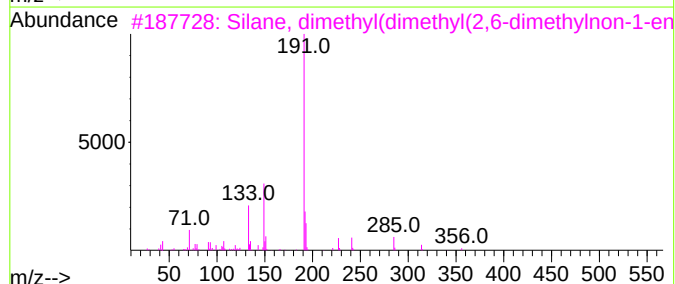
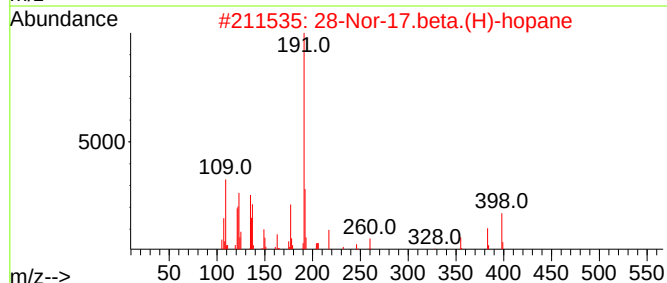
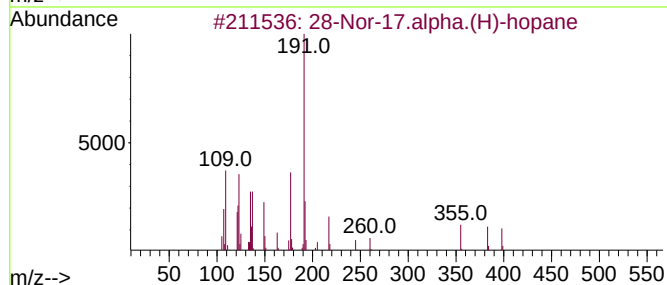
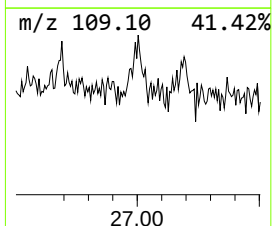
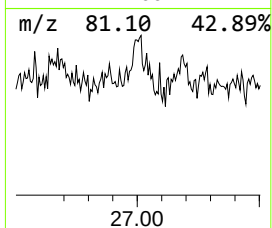
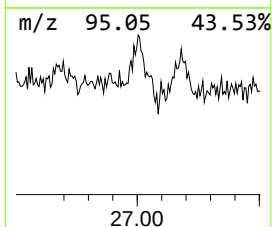
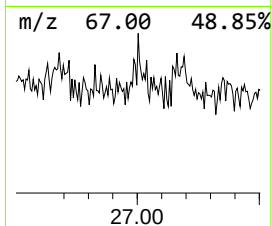
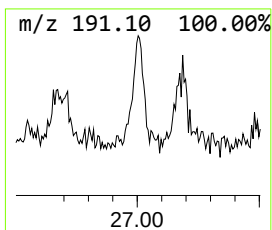
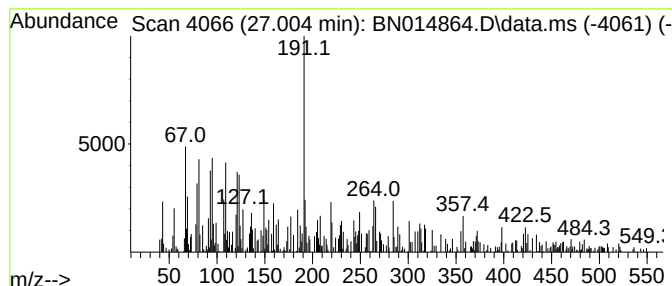
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 unknown-06 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.004	3.19 ng/ul	513486	Perylene-d12	23.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	28-Nor-17.alpha.(H)-hopane			398	C29H50	053584-60-4	49
2	28-Nor-17.beta.(H)-hopane			398	C29H50	036728-72-0	45
3	Silane, dimethyl(dimethyl(2,6-di...			356	C18H36O3Si2	1000347-91-3	27
4	(1,3-Diphenylprop-2-ynyl)diethyl...			263	C19H21N	1000188-31-2	27
5	5,5-Dimethyl-2-(2,2,2-trifluoro-...			249	C8H9F6NO	020967-43-5	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
 Data File : BN014864.D
 Acq On : 10 Jun 2021 14:36
 Operator : CG/JU
 Sample : M2618-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BG5R6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN052521.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-(2-Methyl-...	22.486	2.2	ng/ul	348885	6	23.569	3218260	20.0
unknown-01	22.798	2.5	ng/ul	403076	6	23.569	3218260	20.0
unknown-02	22.945	2.1	ng/ul	339689	6	23.569	3218260	20.0
unknown-03	23.080	5.2	ng/ul	841357	6	23.569	3218260	20.0
5-(1-Isopropeny...	23.780	5.7	ng/ul	912579	6	23.569	3218260	20.0
unknown-04	23.880	2.4	ng/ul	384920	6	23.569	3218260	20.0
Silane, diethyl...	23.927	2.1	ng/ul	332816	6	23.569	3218260	20.0
unknown-05	24.116	2.6	ng/ul	418320	6	23.569	3218260	20.0
Cholestan-3-one	24.339	12.1	ng/ul	1938560	6	23.569	3218260	20.0
28-Nor-17.alpha...	24.645	11.6	ng/ul	1872660	6	23.569	3218260	20.0
(DEL) Alkane: S...	24.886	3.3	ng/ul	528969	6	23.569	3218260	20.0
2,3(1H)-Dihydro...	25.098	3.1	ng/ul	505084	6	23.569	3218260	20.0
2H-1,2,3-Triazo...	25.327	10.7	ng/ul	1720490	6	23.569	3218260	20.0
(DEL) Alkane: S...	25.769	5.5	ng/ul	884435	6	23.569	3218260	20.0
Phenanthrene, 2...	26.210	10.3	ng/ul	1652930	6	23.569	3218260	20.0
unknown-06	27.004	3.2	ng/ul	513486	6	23.569	3218260	20.0