

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050523\
 Data File : BP014979.D
 Acq On : 05 May 2023 12:35
 Operator : CG/JU
 Sample : 02479-16 5X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EW975

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP014979.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.011	102	106	114	rVB	57121	83087	0.58%	0.095%
2	4.246	142	146	153	rVB	24273	32794	0.23%	0.038%
3	4.475	182	185	191	rVB	27417	33115	0.23%	0.038%
4	5.299	320	325	331	rVB2	23070	34977	0.24%	0.040%
5	6.111	458	463	471	rBV	28734	47073	0.33%	0.054%
6	7.311	659	667	676	rBV	169564	331906	2.31%	0.381%
7	7.481	690	696	703	rBV	137160	217344	1.51%	0.250%
8	7.687	725	731	739	rVB	172625	275056	1.92%	0.316%
9	8.163	801	812	820	rBV	1172247	1900080	13.23%	2.183%
10	8.240	820	825	831	rVB	19418	33137	0.23%	0.038%
11	8.410	848	854	867	rBV6	14011	36559	0.25%	0.042%
12	8.875	927	933	942	rBV	192764	322388	2.25%	0.370%
13	8.999	948	954	962	rVB	50438	90568	0.63%	0.104%
14	9.281	994	1002	1006	rBV	23325	43361	0.30%	0.050%
15	9.340	1006	1012	1019	rVV	157697	266986	1.86%	0.307%
16	10.069	1130	1136	1147	rVB2	128776	222348	1.55%	0.255%
17	10.199	1147	1158	1161	rBV8	11562	28224	0.20%	0.032%
18	10.616	1219	1229	1236	rBV	208055	368799	2.57%	0.424%
19	10.999	1283	1294	1300	rBV	1746836	2999755	20.89%	3.447%
20	11.052	1300	1303	1310	rVV	79598	132984	0.93%	0.153%
21	11.140	1311	1318	1327	rVV	72039	141392	0.98%	0.162%
22	12.004	1456	1465	1470	rBV6	16665	36344	0.25%	0.042%
23	12.399	1525	1532	1537	rBV4	21851	37783	0.26%	0.043%
24	12.646	1565	1574	1581	rVB	163474	285777	1.99%	0.328%
25	12.863	1605	1611	1619	rVB	84469	138025	0.96%	0.159%
26	13.840	1773	1777	1788	rVB	31278	65888	0.46%	0.076%
27	13.969	1788	1799	1801	rBV2	82091	134857	0.94%	0.155%
28	14.134	1820	1827	1832	rBV3	135205	253105	1.76%	0.291%
29	14.204	1832	1839	1848	rVV	426389	786007	5.47%	0.903%
30	14.275	1848	1851	1855	rVB2	23572	31723	0.22%	0.036%
31	14.363	1862	1866	1870	rVV	42752	64587	0.45%	0.074%
32	14.504	1884	1890	1901	rBV	523831	842639	5.87%	0.968%
33	14.616	1906	1909	1916	rVB	39294	59821	0.42%	0.069%
34	14.687	1918	1921	1923	rBV	24893	28900	0.20%	0.033%
35	14.716	1923	1926	1930	rVB3	34560	45194	0.31%	0.052%
36	14.810	1931	1942	1948	rBV	2823025	4154844	28.94%	4.774%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
 Title : SVOA CALIBRATION

37	14.875	1948	1953	1960	rVB4	66407	173372	1.21%	0.199%
38	14.998	1968	1974	1983	rBV3	115819	245402	1.71%	0.282%
39	15.204	2003	2009	2015	rBV2	101186	166348	1.16%	0.191%
40	15.281	2018	2022	2029	rVB	52950	74240	0.52%	0.085%
41	15.422	2041	2046	2050	rBV2	35275	63974	0.45%	0.074%
42	15.475	2050	2055	2062	rVV	43859	67549	0.47%	0.078%
43	15.592	2069	2075	2078	rBV4	44382	79244	0.55%	0.091%
44	15.634	2078	2082	2086	rVB2	51739	78588	0.55%	0.090%
45	15.798	2105	2110	2115	rBV	674337	926004	6.45%	1.064%
46	15.851	2115	2119	2123	rVV2	61192	96815	0.67%	0.111%
47	15.916	2125	2130	2136	rVB2	139763	207320	1.44%	0.238%
48	16.157	2166	2171	2176	rVB2	113036	175598	1.22%	0.202%
49	16.298	2192	2195	2199	rVV	30420	43687	0.30%	0.050%
50	16.345	2199	2203	2209	rVV4	53178	91277	0.64%	0.105%
51	16.539	2226	2236	2247	rBV	2558853	3423060	23.84%	3.933%
52	16.839	2282	2287	2288	rBV3	20494	33044	0.23%	0.038%
53	17.092	2327	2330	2335	rBV6	22636	38794	0.27%	0.045%
54	17.216	2347	2351	2354	rBV	59657	85967	0.60%	0.099%
55	17.251	2354	2357	2360	rVV2	46765	62184	0.43%	0.071%
56	17.304	2362	2366	2370	rVV5	74309	138091	0.96%	0.159%
57	17.351	2371	2374	2377	rVV2	51093	80282	0.56%	0.092%
58	17.381	2377	2379	2384	rVV	59711	79603	0.55%	0.091%
59	17.563	2404	2410	2415	rBV	3301688	5001198	34.83%	5.746%
60	17.610	2415	2418	2422	rVV	988007	1274738	8.88%	1.465%
61	17.663	2422	2427	2431	rVV	736444	1054416	7.34%	1.211%
62	17.698	2431	2433	2437	rVB	192396	202682	1.41%	0.233%
63	17.963	2474	2478	2481	rBV	86936	128989	0.90%	0.148%
64	18.034	2487	2490	2494	rVB3	52270	62692	0.44%	0.072%
65	18.139	2504	2508	2515	rVB5	110868	221008	1.54%	0.254%
66	18.422	2551	2556	2560	rBV	381977	547739	3.82%	0.629%
67	18.469	2560	2564	2573	rVB2	217024	342078	2.38%	0.393%
68	18.604	2582	2587	2591	rBV2	235069	425561	2.96%	0.489%
69	18.645	2591	2594	2599	rVB	109296	157330	1.10%	0.181%
70	18.704	2600	2604	2609	rBV4	65167	111622	0.78%	0.128%
71	18.833	2623	2626	2630	rBV2	171183	231851	1.61%	0.266%
72	18.898	2634	2637	2640	rVV	92535	116143	0.81%	0.133%
73	18.939	2641	2644	2648	rVB3	77486	99786	0.70%	0.115%
74	19.245	2693	2696	2701	rVB	286235	364374	2.54%	0.419%
75	19.298	2702	2705	2709	rVB2	122172	165222	1.15%	0.190%
76	19.386	2718	2720	2723	rBV	64669	61445	0.43%	0.071%

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77	19.557	2744	2749	2756	rVB	2056566	2623386	18.27%	3.014%
78	19.645	2761	2764	2768	rVB4	142357	177618	1.24%	0.204%
79	19.880	2799	2804	2806	rBV2	1141903	1581076	11.01%	1.817%
80	19.910	2806	2809	2816	rVB	1714903	2225363	15.50%	2.557%
81	20.216	2857	2861	2867	rBV4	178089	326233	2.27%	0.375%
82	20.380	2886	2889	2893	rVB2	222819	322042	2.24%	0.370%
83	20.480	2904	2906	2909	rBV	92164	91582	0.64%	0.105%
84	20.680	2936	2940	2944	rBV	181115	261225	1.82%	0.300%
85	20.763	2951	2954	2957	rVB	173853	189432	1.32%	0.218%
86	20.863	2968	2971	2977	rBV2	213864	256556	1.79%	0.295%
87	21.169	3020	3023	3026	rVB	228840	229396	1.60%	0.264%
88	21.316	3045	3048	3052	rVB	485648	533111	3.71%	0.613%
89	21.545	3079	3087	3096	rBV2	8270973	14357092	100.00%	16.495%
90	21.645	3097	3104	3108	rVV2	3380967	5905642	41.13%	6.785%
91	21.686	3108	3111	3120	rVB	735602	1027176	7.15%	1.180%
92	22.057	3170	3174	3185	rVB	1394511	2130802	14.84%	2.448%
93	22.263	3206	3209	3212	rBV2	326204	467000	3.25%	0.537%
94	22.298	3212	3215	3217	rVV	609997	819683	5.71%	0.942%
95	22.345	3217	3223	3235	rVB	6651630	10434903	72.68%	11.989%
96	22.616	3265	3269	3281	rVB	1511448	2732232	19.03%	3.139%
97	22.769	3292	3295	3299	rBV	374496	568921	3.96%	0.654%
98	22.827	3300	3305	3317	rVB	1534664	2722523	18.96%	3.128%
99	23.145	3355	3359	3368	rVB	1152448	2081535	14.50%	2.392%
100	24.245	3540	3546	3553	rBV2	1522651	3365338	23.44%	3.867%

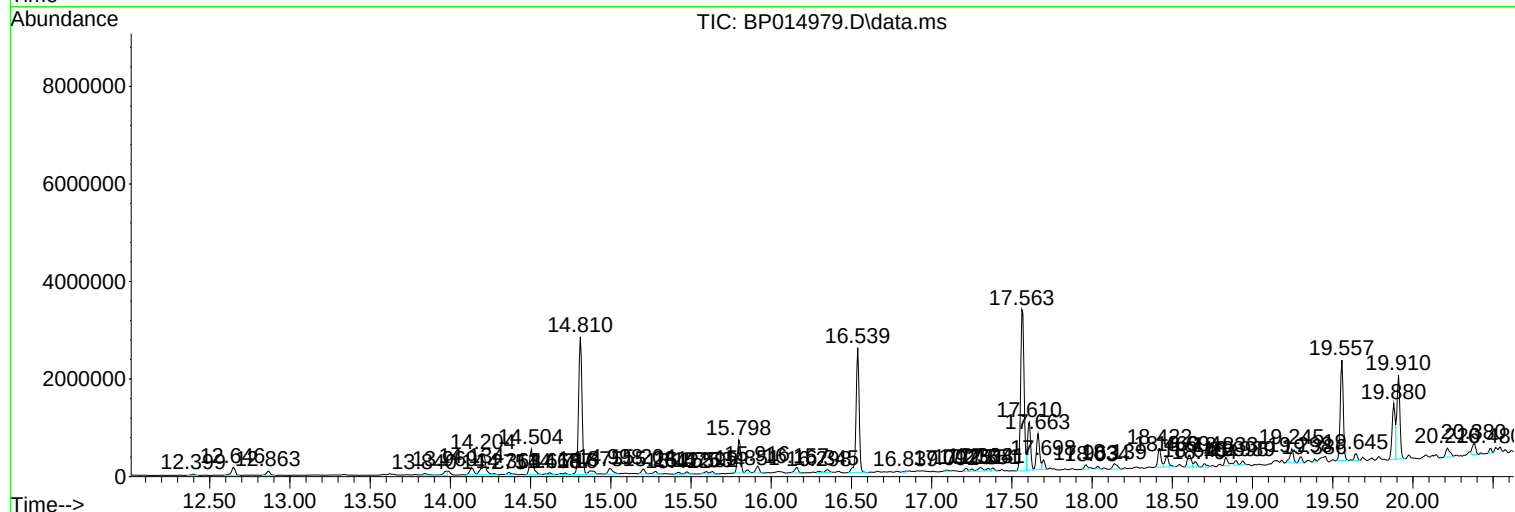
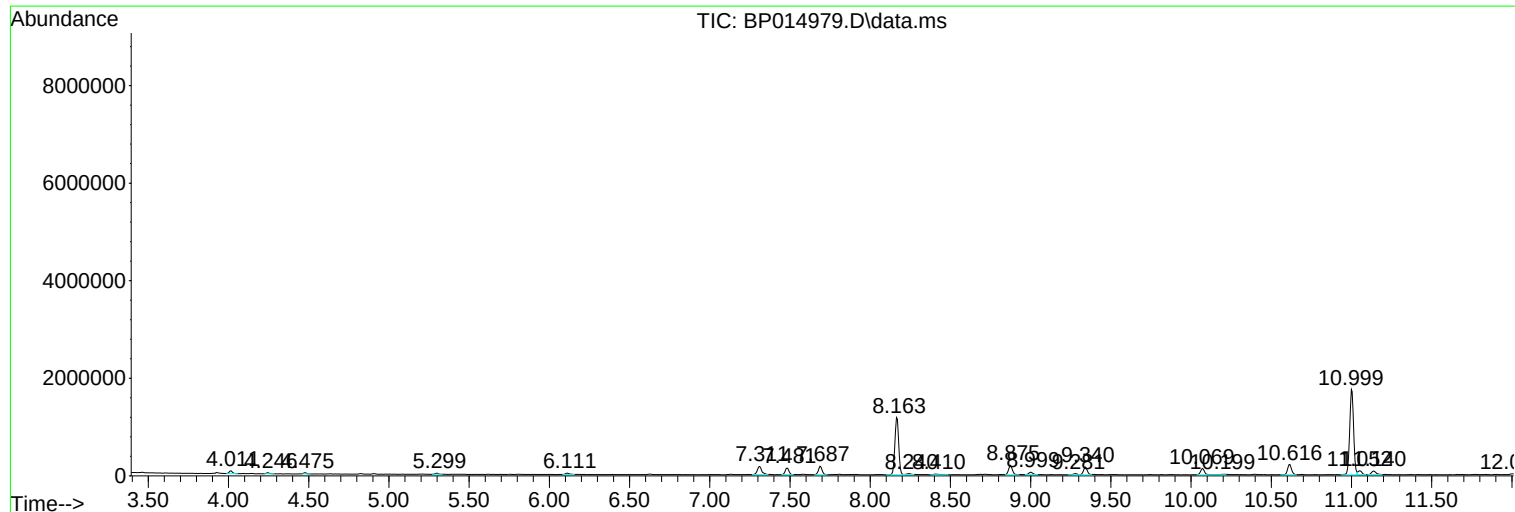
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



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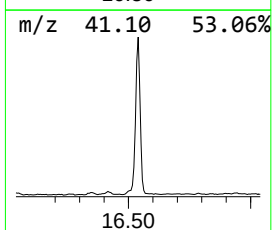
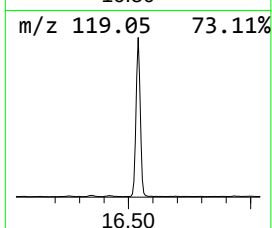
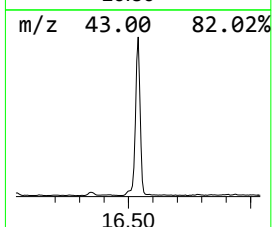
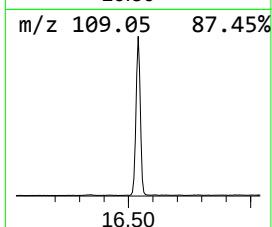
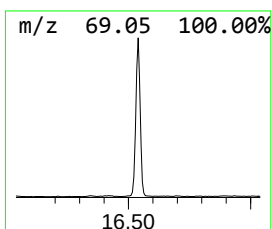
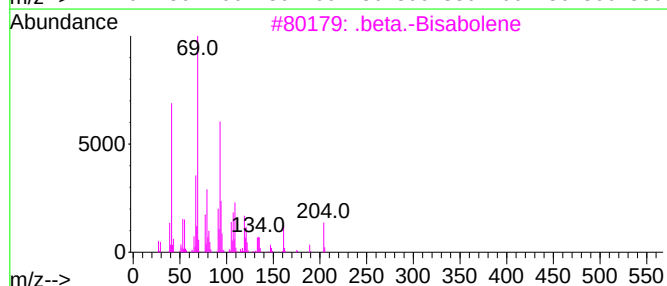
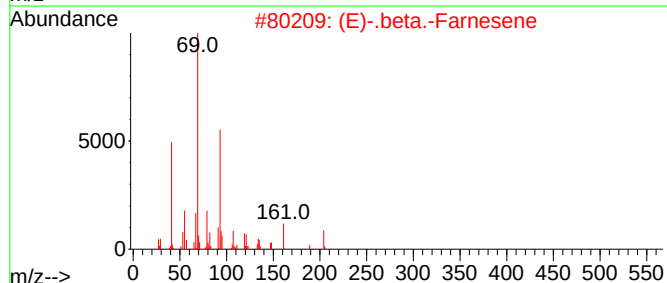
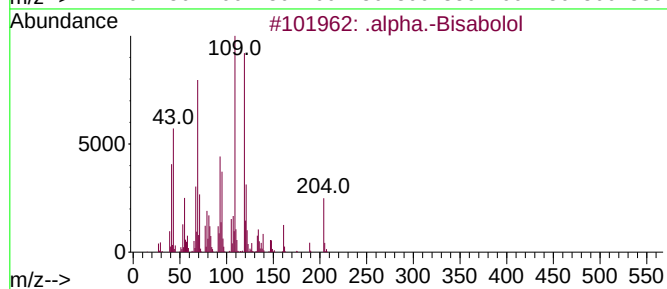
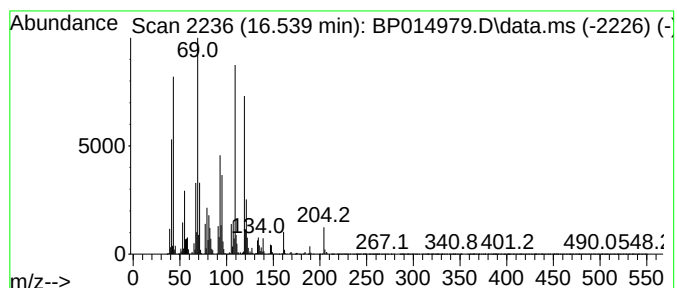
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 .alpha.-Bisabolol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.540	13.69 ng/ul	3423060	Phenanthrene-d10	17.563	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	.alpha.-Bisabolol	222	C15H26O	000515-69-5	91
2	(E)-.beta.-Farnesene	204	C15H24	018794-84-8	87
3	.beta.-Bisabolene	204	C15H24	000495-61-4	70
4	(E)-.beta.-Farnesene	204	C15H24	018794-84-8	55
5	Tricyclo[5.4.0.0(2,8)]undec-9-en...	204	C15H24	005989-08-2	47



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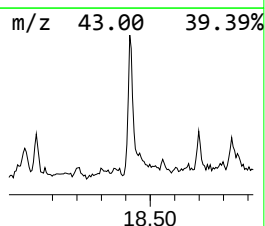
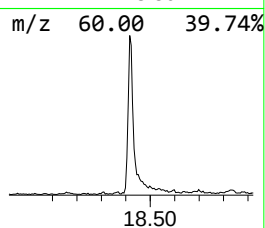
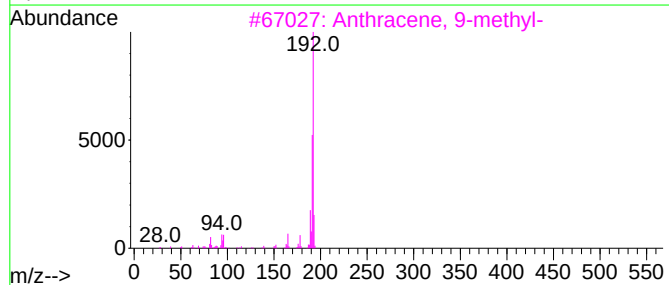
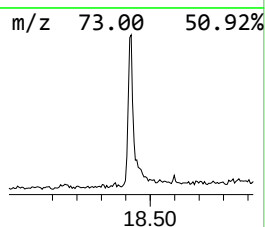
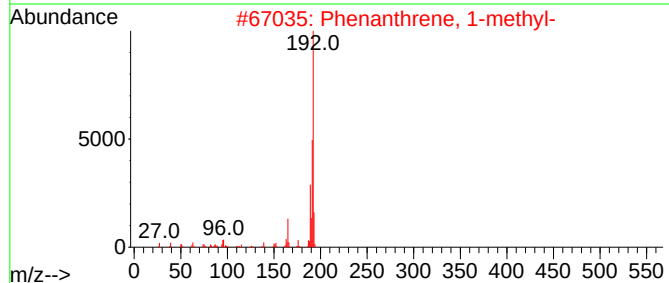
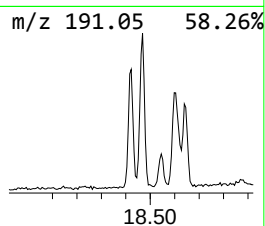
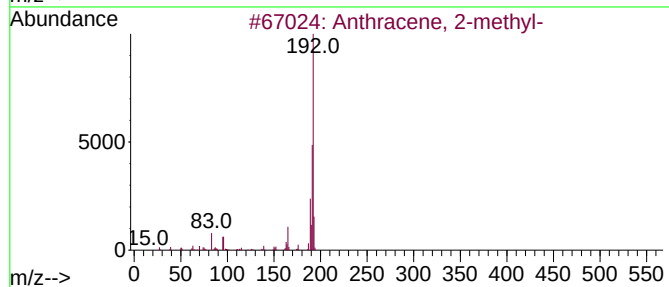
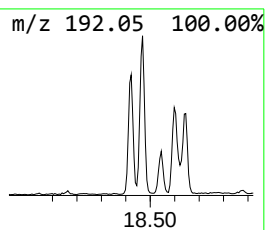
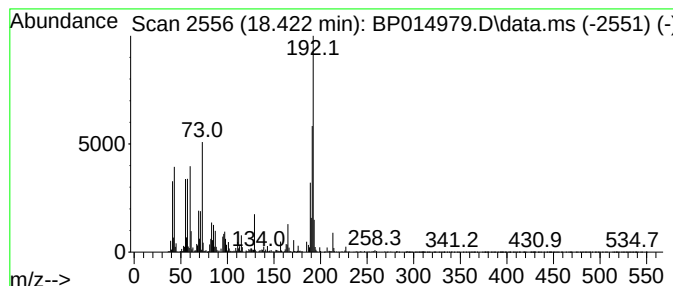
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Anthracene, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.422	2.19 ng/ul	547739	Phenanthrene-d10	17.563

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Anthracene, 2-methyl-	192	C15H12	000613-12-7	92
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	91
3			Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
4			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	90
5			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	78



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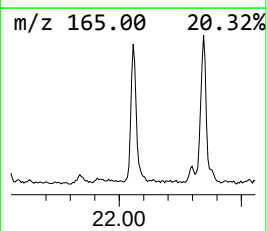
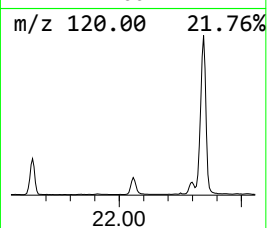
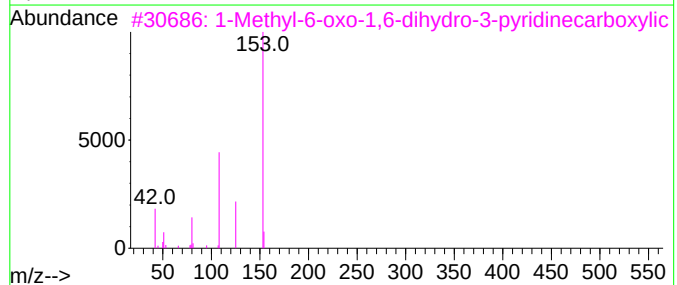
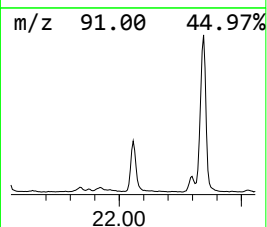
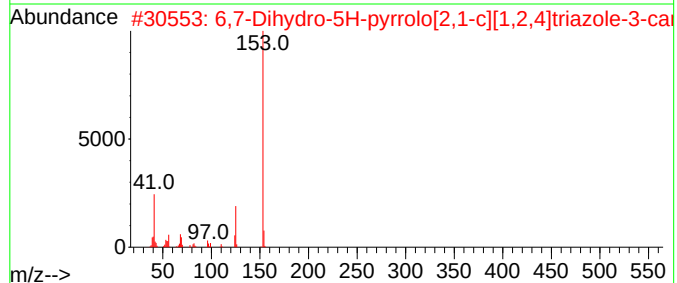
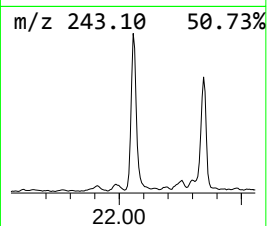
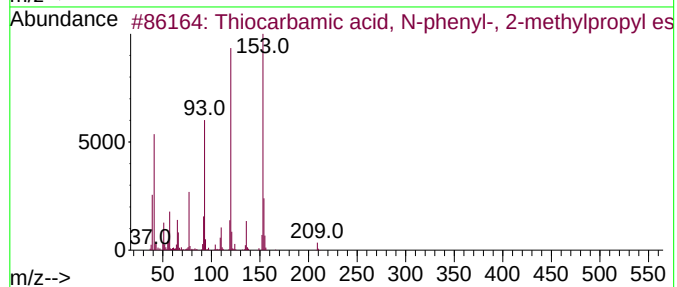
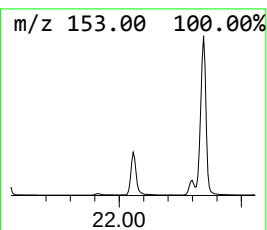
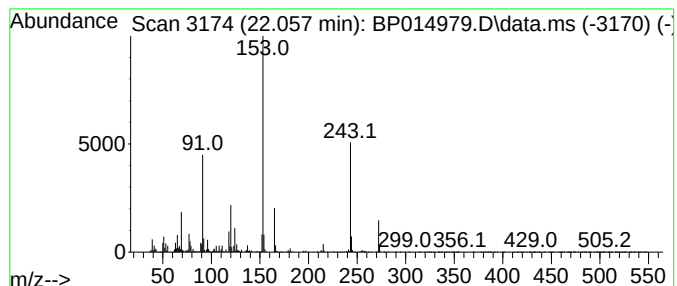
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.057	7.22 ng/ul	2130800	Chrysene-d12	21.645

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Thiocarbamic acid, N-phenyl-, 2-...	209	C11H15NOS	027830-90-6	32
2			6,7-Dihydro-5H-pyrrolo[2,1-c][1,...	153	C6H7N3O2	1000296-76-1	27
3			1-Methyl-6-oxo-1,6-dihydro-3-pyr...	153	C7H7N03	003719-45-7	27
4			2-Methyl-1-(2,4,6-trihydroxyphen...	196	C10H12O4	035458-21-0	27
5			Preleptospermone	210	C11H14O4	026103-97-9	27



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Data File : BP014979.D
Acq On : 05 May 2023 12:35
Operator : CG/JU
Sample : 02479-16 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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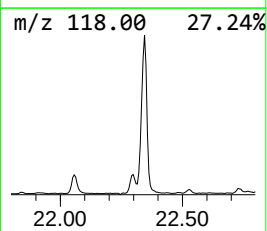
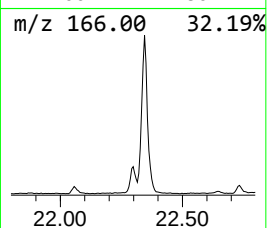
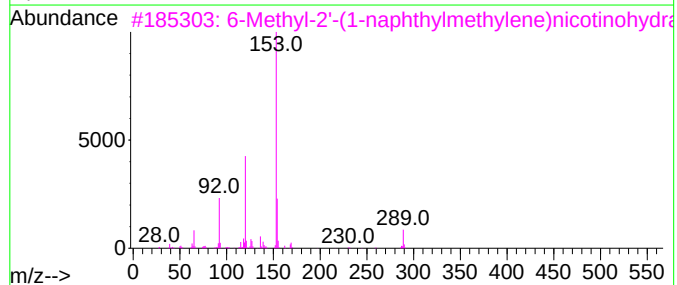
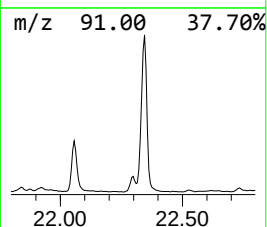
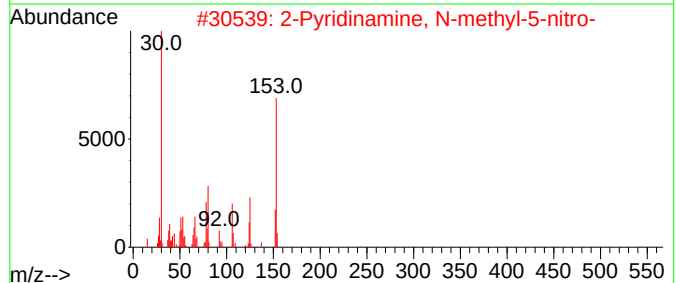
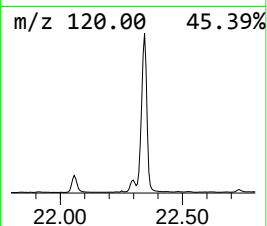
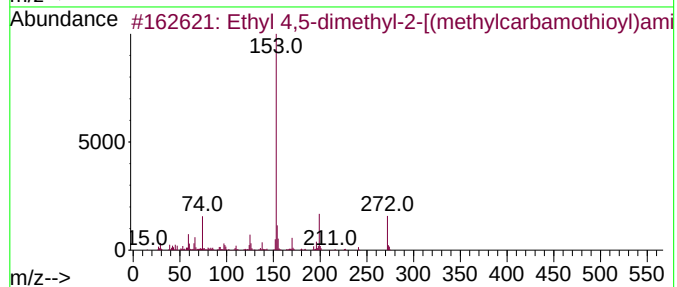
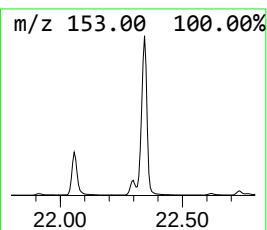
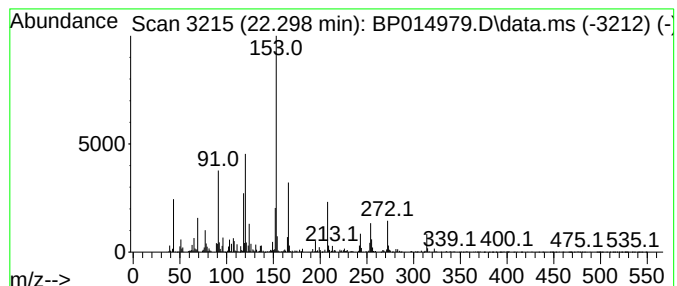
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown-02 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.298	2.78 ng/ul	819683	Chrysene-d12	21.645

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethyl 4,5-dimethyl-2-[(methylcar...	272	C11H16N2O2S2	059898-39-4	27
2			2-Pyridinamine, N-methyl-5-nitro-	153	C6H7N3O2	004093-89-4	25
3			6-Methyl-2'-(1-naphthylmethylene...	289	C18H15N3O	1000239-87-7	25
4			Naringenin	272	C15H12O5	000480-41-1	18
5			N-(4-Chloro-2-fluorophenyl)-2-(2...	297	C12H9ClFN3OS	317325-22-7	17



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050523\
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Sample : 02479-16 5X
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ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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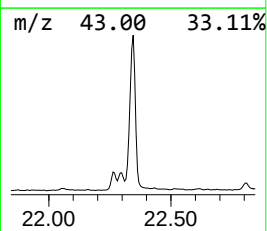
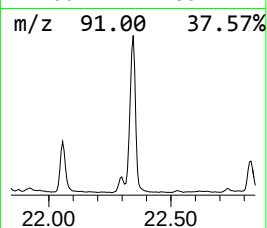
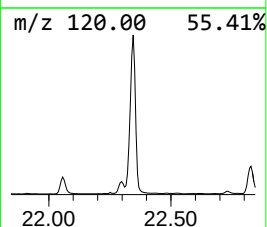
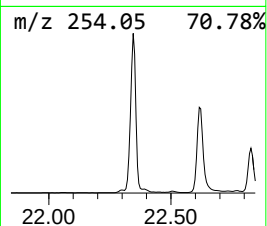
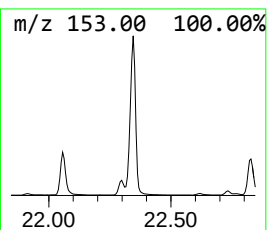
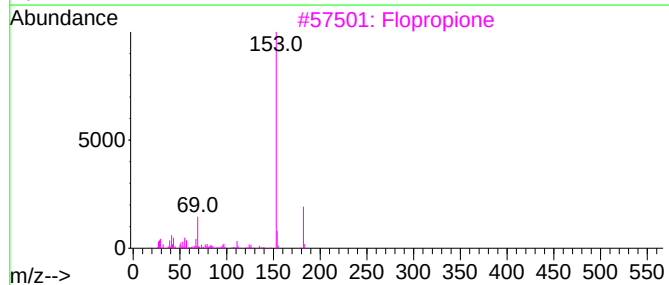
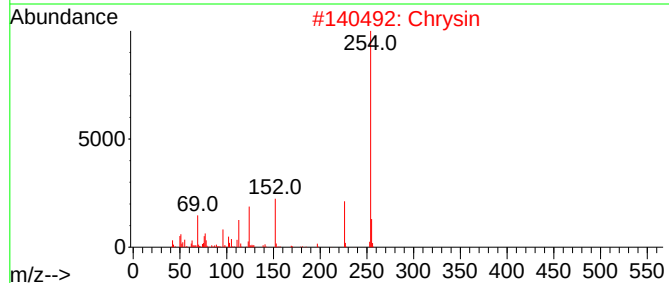
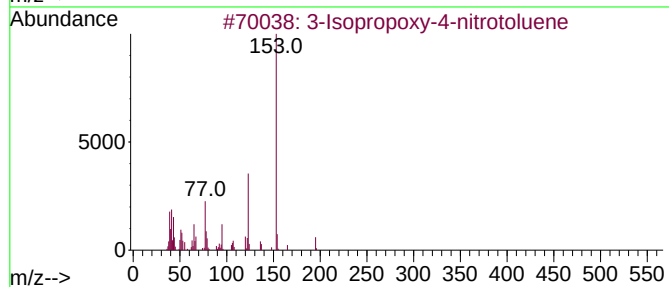
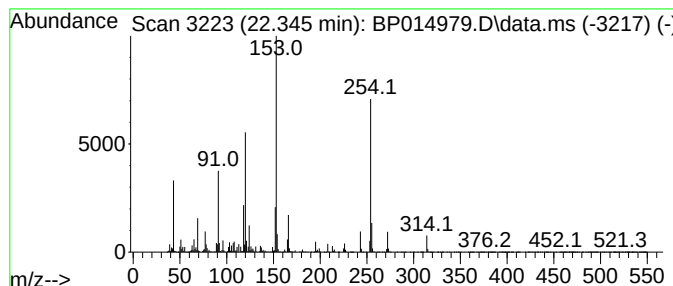
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown-03 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.345	35.34 ng/ul	10434900	Chrysene-d12	21.645

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Isopropoxy-4-nitrotoluene	195	C10H13NO3	126802-98-0	22
2			Chrysin	254	C15H10O4	000480-40-0	15
3			Flopropione	182	C9H10O4	002295-58-1	14
4			1,3-Benzenediol, 5-(dimethylamino)-	153	C8H11NO2	040248-00-8	14
5			1-Methyl-6-oxo-1,6-dihydro-3-pyr...	153	C7H7NO3	003719-45-7	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050523\
Data File : BP014979.D
Acq On : 05 May 2023 12:35
Operator : CG/JU
Sample : 02479-16 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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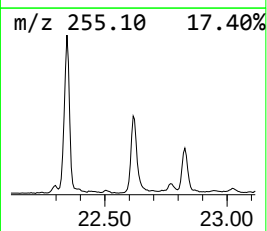
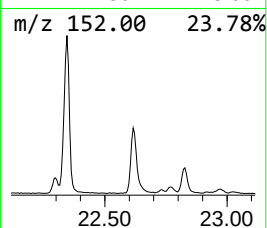
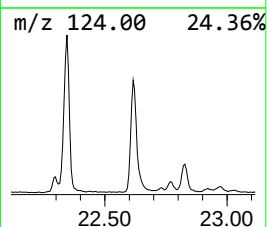
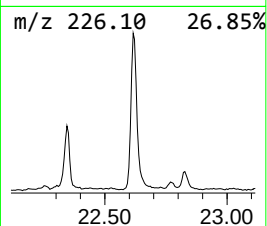
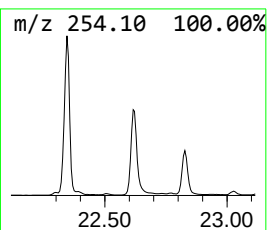
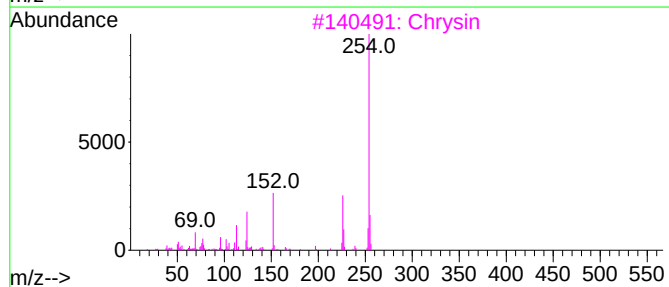
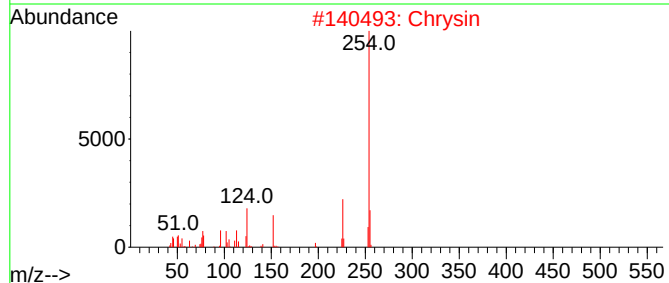
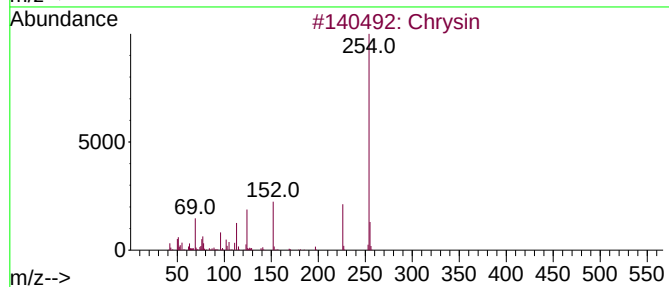
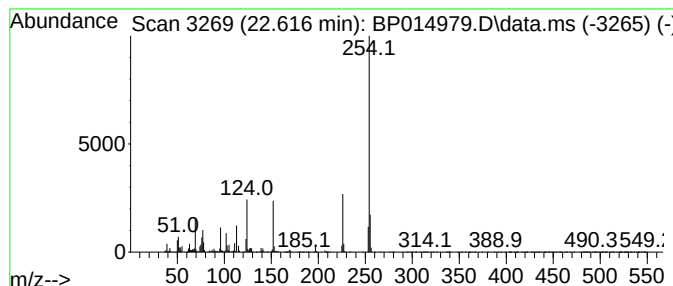
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Chrysin Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.616	9.25 ng/ul	2732230	Chrysene-d12	21.645

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Chrysin	254	C15H10O4	000480-40-0	98
2			Chrysin	254	C15H10O4	000480-40-0	95
3			Chrysin	254	C15H10O4	000480-40-0	94
4			5,6-Dihydroxyflavone	254	C15H10O4	006665-66-3	68
5			2,2'-Binaphthalene	254	C20H14	000612-78-2	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050523\
Data File : BP014979.D
Acq On : 05 May 2023 12:35
Operator : CG/JU
Sample : 02479-16 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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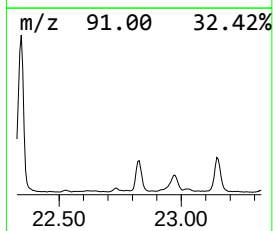
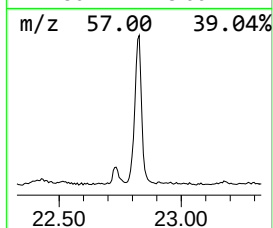
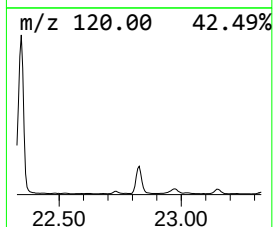
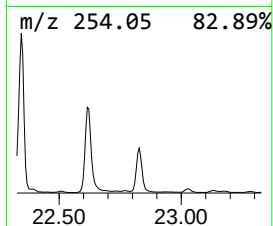
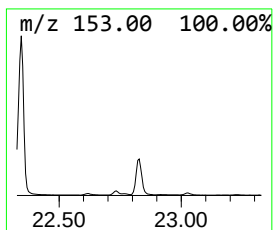
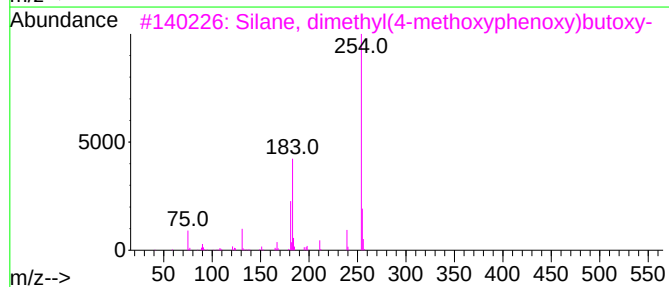
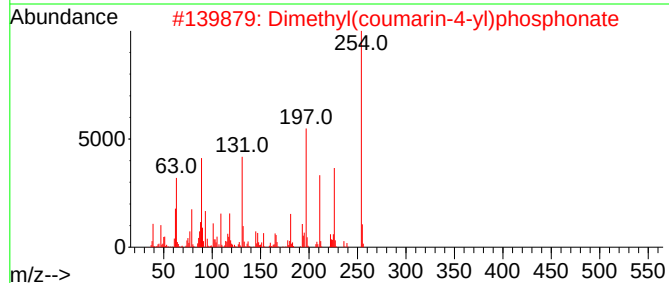
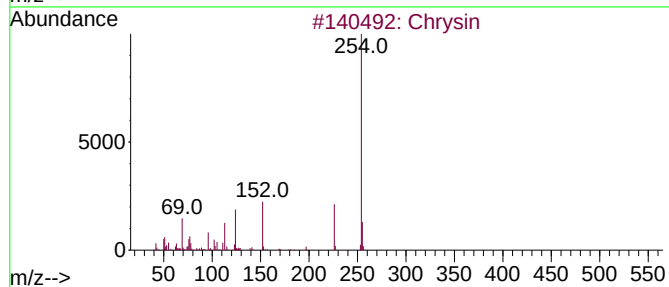
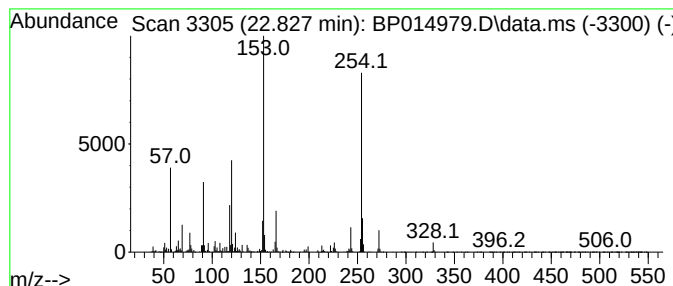
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 unknown-04 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.827	9.22 ng/ul	2722520	Chrysene-d12	21.645

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Chrysin	254	C15H10O4	000480-40-0	22
2			Dimethyl(coumarin-4-yl)phosphonate	254	C11H11O5P	139300-89-3	22
3			Silane, dimethyl(4-methoxyphenox...	254	C13H22O3Si	1000347-14-2	14
4			Naphthalen-1(2H)-one, 3,4-dihydr...	254	C16H14O5	351066-46-1	14
5			1,1'-Binaphthalene	254	C20H14	000604-53-5	14



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050523\
Data File : BP014979.D
Acq On : 05 May 2023 12:35
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Misc :
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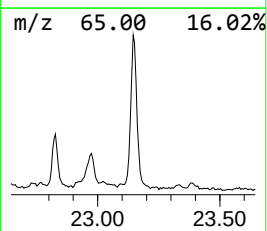
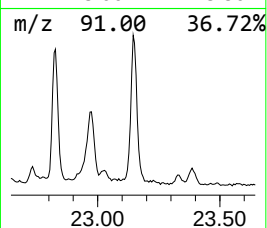
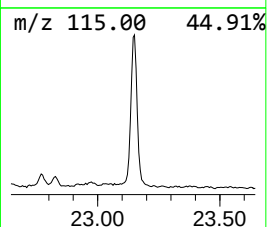
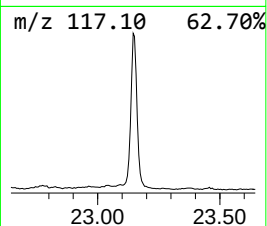
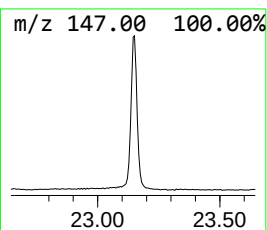
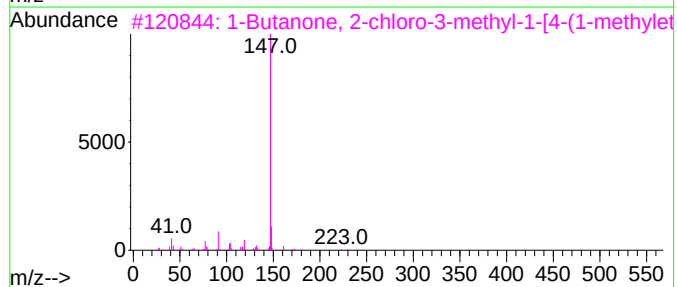
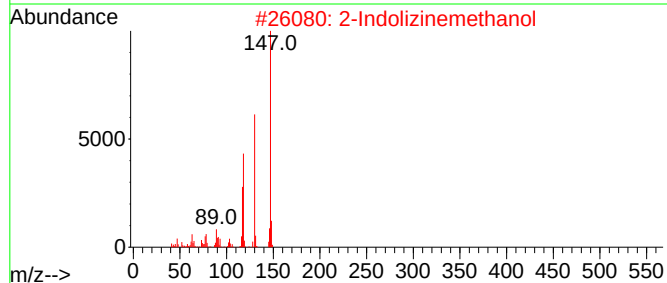
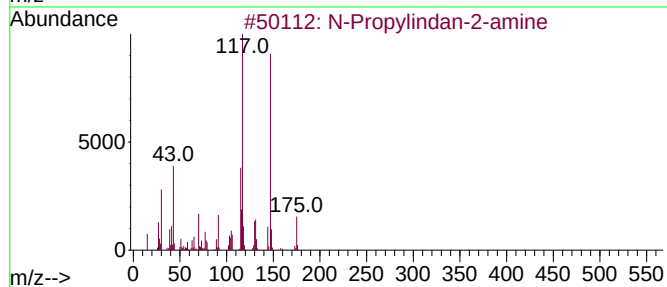
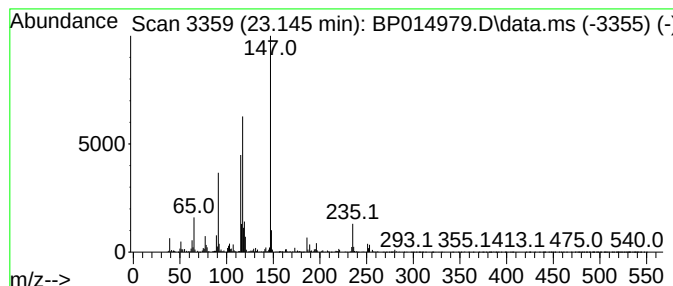
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 N-Propylindan-2-amine Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.145	12.37 ng/ul	2081540	Perylene-d12	24.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			N-Propylindan-2-amine	175	C12H17N	082985-09-9	59
2			2-Indolizinemethanol	147	C9H9NO	022320-21-4	38
3			1-Butanone, 2-chloro-3-methyl-1-...	238	C14H19ClO	055955-90-3	38
4			Benzoic acid, 4-propyl-, 4-cyano...	265	C17H15NO2	056131-49-8	38
5			1-Naphthalenamine, 5,6,7,8-tetra...	147	C10H13N	002217-41-6	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050523\
Data File : BP014979.D
Acq On : 05 May 2023 12:35
Operator : CG/JU
Sample : 02479-16 5X
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Anthracene, 2-m...	18.422	2.2	ng/ul	547739	4	17.563	5001200	20.0
unknown-01	22.057	7.2	ng/ul	2130800	5	21.645	5905640	20.0
unknown-02	22.298	2.8	ng/ul	819683	5	21.645	5905640	20.0
unknown-03	22.345	35.3	ng/ul	10434900	5	21.645	5905640	20.0
Chrysin	22.616	9.3	ng/ul	2732230	5	21.645	5905640	20.0
unknown-04	22.827	9.2	ng/ul	2722520	5	21.645	5905640	20.0
N-Propylindan-2...	23.145	12.4	ng/ul	2081540	6	24.245	3365340	20.0