

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112723\
 Data File : BP018452.D
 Acq On : 28 Nov 2023 22:04
 Operator : MA/JU
 Sample : 05332-16DL 5X
 Misc :
 ALS Vial : 61 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YCG97DL

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-BP112223.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP018452.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.687	99	102	111	rBV	80085	132291	1.58%	0.200%
2	7.016	663	668	679	rVB	69631	139325	1.66%	0.211%
3	7.187	691	697	704	rBV	317308	492967	5.87%	0.746%
4	7.381	720	730	738	rBV	270596	446599	5.32%	0.676%
5	7.852	802	810	818	rBV	1330701	2039743	24.31%	3.086%
6	8.228	868	874	880	rBV	40418	63111	0.75%	0.095%
7	8.563	924	931	942	rVB	193244	318386	3.79%	0.482%
8	8.963	993	999	1002	rBV	117165	180407	2.15%	0.273%
9	9.010	1002	1007	1019	rVB	282001	509582	6.07%	0.771%
10	9.293	1049	1055	1061	rBV2	27707	57549	0.69%	0.087%
11	9.487	1083	1088	1093	rVB	73221	112812	1.34%	0.171%
12	9.734	1122	1130	1137	rBV	164548	278525	3.32%	0.421%
13	10.051	1177	1184	1193	rBV4	93658	189407	2.26%	0.287%
14	10.269	1213	1221	1230	rBV	318222	550220	6.56%	0.833%
15	10.651	1274	1286	1292	rVV	1711267	2817197	33.57%	4.263%
16	10.704	1292	1295	1301	rVV3	40335	68530	0.82%	0.104%
17	10.793	1301	1310	1314	rVV	252471	488389	5.82%	0.739%
18	10.840	1314	1318	1325	rVV2	286611	548126	6.53%	0.829%
19	10.898	1325	1328	1332	rVV	67049	113901	1.36%	0.172%
20	10.951	1332	1337	1343	rVV2	88004	161565	1.93%	0.244%
21	11.022	1345	1349	1353	rVV2	35319	60957	0.73%	0.092%
22	11.075	1353	1358	1364	rVV	160509	303282	3.61%	0.459%
23	11.134	1364	1368	1373	rVB2	38356	61787	0.74%	0.093%
24	11.240	1380	1386	1392	rBV	37434	61644	0.73%	0.093%
25	11.316	1392	1399	1405	rVB2	37933	71013	0.85%	0.107%
26	11.422	1410	1417	1422	rBV	54377	89882	1.07%	0.136%
27	11.569	1436	1442	1446	rBV	38599	70265	0.84%	0.106%
28	11.698	1460	1464	1468	rVV2	57350	97558	1.16%	0.148%
29	11.757	1468	1474	1478	rVV4	46043	113228	1.35%	0.171%
30	11.846	1484	1489	1498	rVV2	74849	138533	1.65%	0.210%
31	12.040	1517	1522	1529	rVB2	46252	73455	0.88%	0.111%
32	12.151	1535	1541	1546	rBV2	28698	75754	0.90%	0.115%
33	12.210	1546	1551	1561	rVV	137950	262241	3.13%	0.397%
34	12.375	1574	1579	1583	rBV	71264	114522	1.36%	0.173%
35	12.428	1583	1588	1594	rVV5	31843	69820	0.83%	0.106%
36	12.498	1594	1600	1611	rVB5	85085	242353	2.89%	0.367%

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Integrator: RTE

Smoothing : OFF

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Stop Thrs : 0

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
 Title : SVOA CALIBRATION

37	13.328	1734	1741	1747	rVV	1755588	2514376	29.96%	3.805%
38	13.392	1747	1752	1759	rVV	79363	147965	1.76%	0.224%
39	13.504	1767	1771	1774	rVV2	35876	55726	0.66%	0.084%
40	13.545	1774	1778	1784	rVB2	78457	119405	1.42%	0.181%
41	13.663	1793	1798	1806	rVV3	35370	80866	0.96%	0.122%
42	13.816	1815	1824	1830	rBV	477776	729046	8.69%	1.103%
43	13.904	1832	1839	1847	rVB	712807	1027799	12.25%	1.555%
44	14.045	1851	1863	1866	rBV	319200	475809	5.67%	0.720%
45	14.081	1866	1869	1874	rVB	177229	236067	2.81%	0.357%
46	14.210	1879	1891	1905	rBV2	5074283	8391670	100.00%	12.698%
47	14.492	1928	1939	1944	rBV	2007399	3169929	37.77%	4.796%
48	14.551	1944	1949	1958	rVB	2753378	4007361	47.75%	6.064%
49	14.687	1968	1972	1982	rBV4	32674	85996	1.02%	0.130%
50	14.887	1996	2006	2015	rBV	1853749	2614063	31.15%	3.955%
51	14.963	2015	2019	2028	rVB2	46243	87314	1.04%	0.132%
52	15.092	2035	2041	2049	rBV2	94883	226300	2.70%	0.342%
53	15.175	2049	2055	2061	rVB2	75798	143860	1.71%	0.218%
54	15.263	2065	2070	2075	rBV2	66433	142832	1.70%	0.216%
55	15.310	2075	2078	2083	rVV2	55325	80300	0.96%	0.122%
56	15.439	2093	2100	2103	rBV	95619	164947	1.97%	0.250%
57	15.481	2103	2107	2112	rVV	1074605	1530527	18.24%	2.316%
58	15.539	2112	2117	2123	rVV	2276068	3221580	38.39%	4.875%
59	15.592	2123	2126	2130	rVV	119816	179745	2.14%	0.272%
60	15.634	2130	2133	2135	rVV	63181	85586	1.02%	0.130%
61	15.663	2135	2138	2141	rVV2	59540	85144	1.01%	0.129%
62	15.698	2141	2144	2149	rVV3	52424	82353	0.98%	0.125%
63	15.751	2149	2153	2156	rVV2	40226	62414	0.74%	0.094%
64	15.786	2156	2159	2163	rVV	97677	130300	1.55%	0.197%
65	15.834	2163	2167	2172	rVB	166198	223978	2.67%	0.339%
66	15.981	2187	2192	2200	rVB	205359	377759	4.50%	0.572%
67	16.069	2204	2207	2215	rVB3	36303	59677	0.71%	0.090%
68	16.216	2227	2232	2238	rBV	109363	162218	1.93%	0.245%
69	16.504	2276	2281	2283	rBV2	43086	62139	0.74%	0.094%
70	16.545	2283	2288	2293	rVV	76280	140861	1.68%	0.213%
71	16.892	2343	2347	2356	rVB	218650	323502	3.86%	0.489%
72	17.063	2372	2376	2382	rVV	150341	205299	2.45%	0.311%
73	17.245	2401	2407	2412	rBV	2474402	3422321	40.78%	5.178%
74	17.286	2412	2414	2418	rVV	227856	279248	3.33%	0.423%
75	17.345	2418	2424	2428	rVV2	1360544	2045023	24.37%	3.094%
76	17.381	2428	2430	2436	rVV	389109	448827	5.35%	0.679%

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77	17.451	2436	2442	2449	rVB	712761	995404	11.86%	1.506%
78	17.651	2466	2476	2481	rBV	591334	886641	10.57%	1.342%
79	17.775	2489	2497	2500	rBV5	34082	96521	1.15%	0.146%
80	17.816	2500	2504	2512	rVB2	63635	128629	1.53%	0.195%
81	18.122	2552	2556	2559	rVV	86605	117189	1.40%	0.177%
82	18.163	2559	2563	2566	rVV2	200270	304496	3.63%	0.461%
83	18.204	2566	2570	2573	rVV	250895	353513	4.21%	0.535%
84	18.239	2573	2576	2582	rVB2	96895	135745	1.62%	0.205%
85	18.310	2582	2588	2592	rBV	259460	424539	5.06%	0.642%
86	18.345	2592	2594	2599	rVB	63397	74454	0.89%	0.113%
87	18.422	2603	2607	2610	rBV3	70096	110542	1.32%	0.167%
88	18.463	2610	2614	2617	rVB	122838	151727	1.81%	0.230%
89	18.551	2624	2629	2636	rBV	194482	286357	3.41%	0.433%
90	18.616	2636	2640	2642	rVV3	68038	95362	1.14%	0.144%
91	18.645	2642	2645	2650	rVB	106762	147755	1.76%	0.224%
92	19.157	2727	2732	2739	rBV3	58261	114503	1.36%	0.173%
93	19.280	2748	2753	2758	rVB	624429	832831	9.92%	1.260%
94	19.363	2764	2767	2773	rVB3	52510	84643	1.01%	0.128%
95	19.610	2803	2809	2812	rBV	1577402	2191923	26.12%	3.317%
96	19.639	2812	2814	2823	rVB	592697	659685	7.86%	0.998%
97	20.022	2876	2879	2884	rVB2	48014	65643	0.78%	0.099%
98	20.086	2885	2890	2896	rBV	234560	369934	4.41%	0.560%
99	21.463	3118	3124	3134	rBV	2548197	4278873	50.99%	6.474%
100	24.251	3589	3598	3616	rVB	892211	3936610	46.91%	5.957%

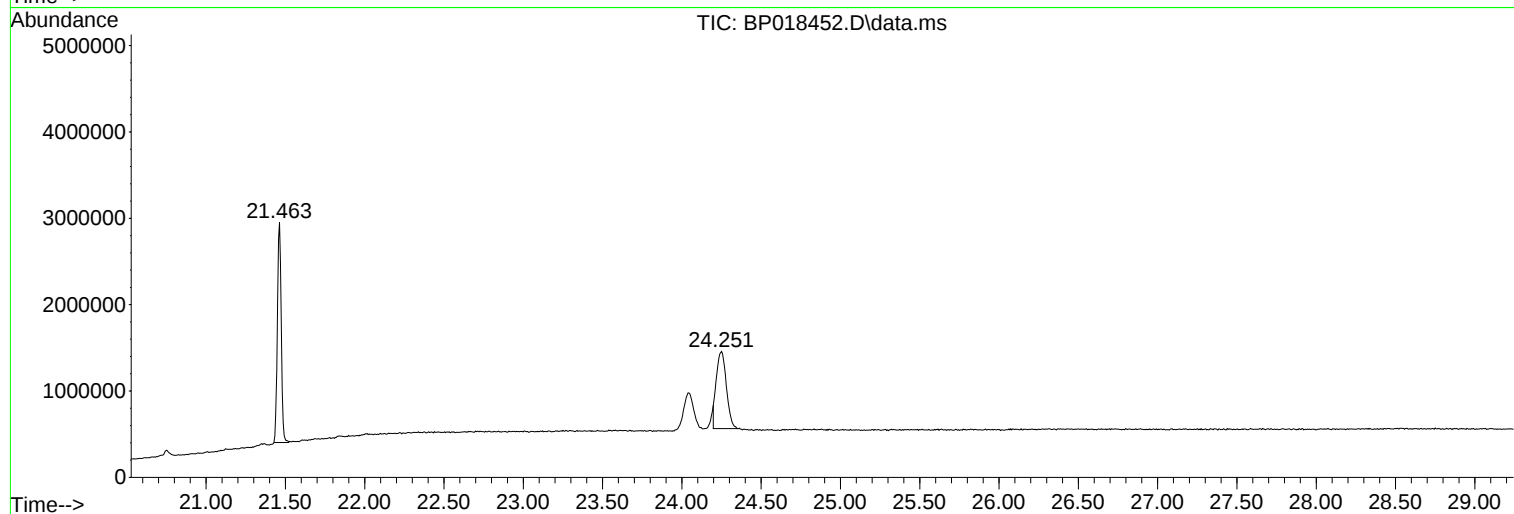
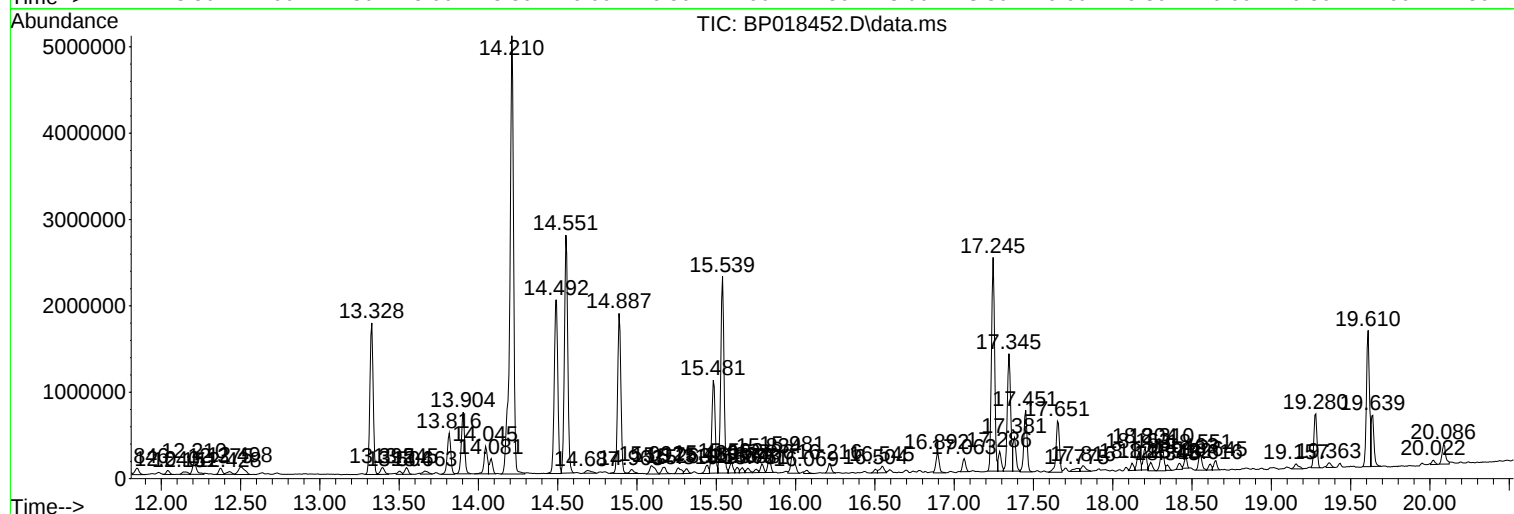
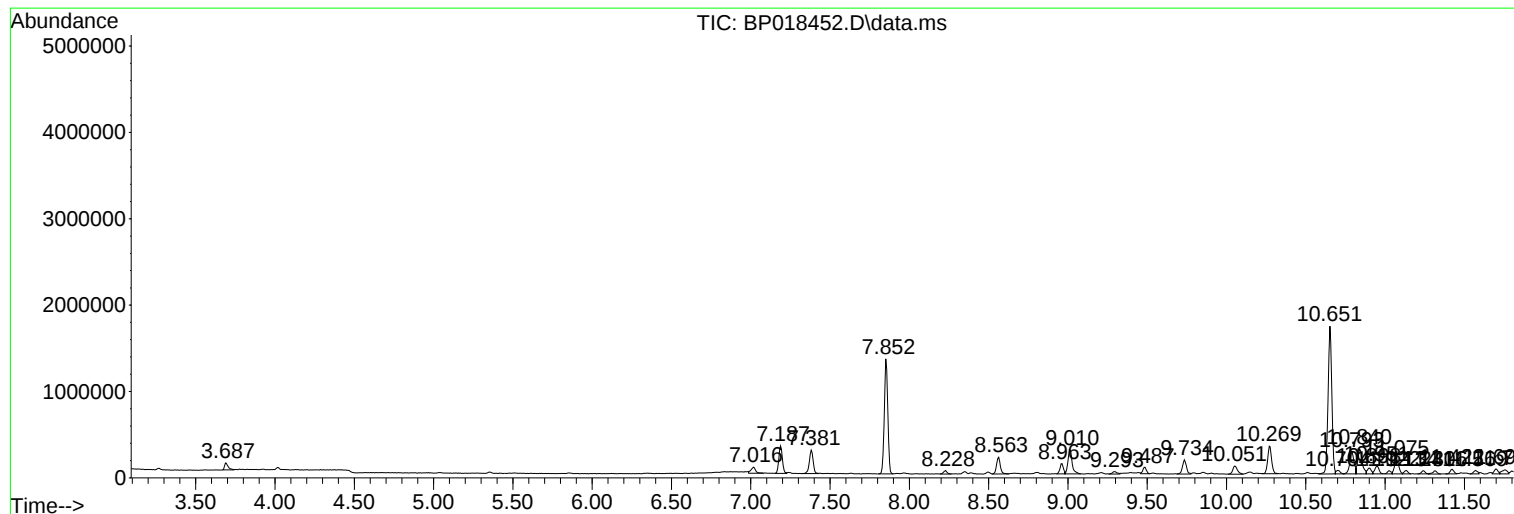
Sum of corrected areas: 66088577

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
Quant Title : SVOA CALIBRATION

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



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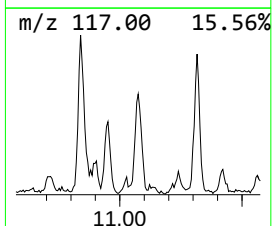
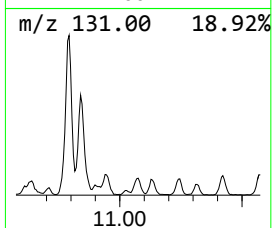
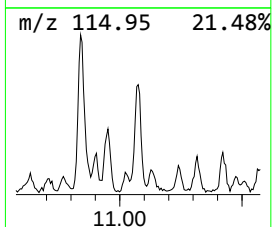
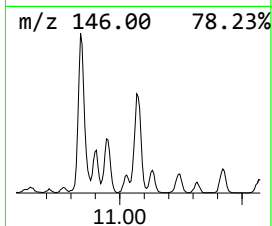
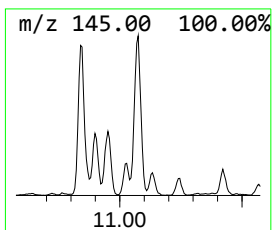
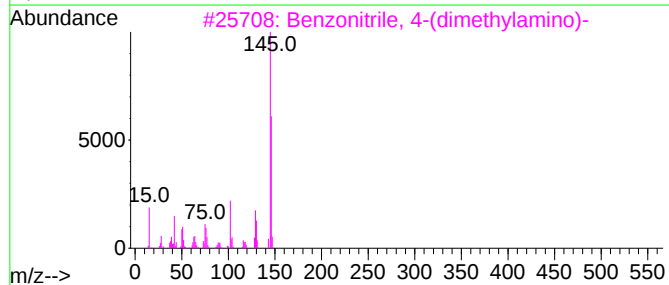
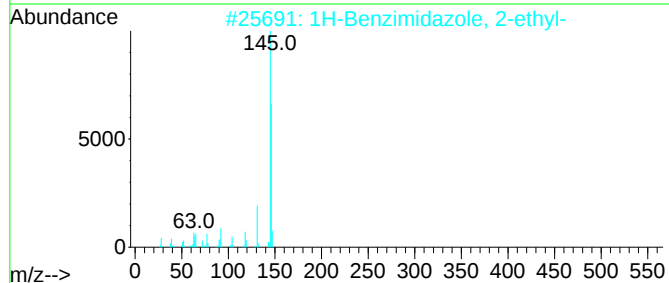
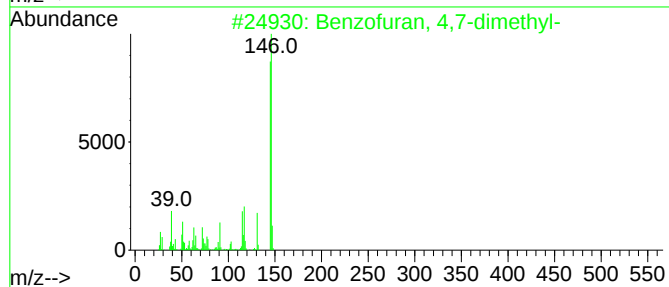
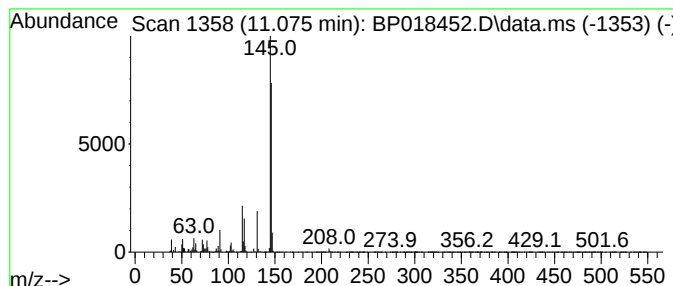
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzofuran, 4,7-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.075	2.15 ng/ul	303282	Naphthalene-d8	10.651

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzofuran, 4,7-dimethyl-	146	C10H10O	028715-26-6	76
2			1H-Benzimidazole, 2-ethyl-	146	C9H10N2	001848-84-6	64
3			Benzonitrile, 4-(dimethylamino)-	146	C9H10N2	001197-19-9	64
4			Benzonitrile, 4-(dimethylamino)-	146	C9H10N2	001197-19-9	59
5			Benzonitrile, 4-(dimethylamino)-	146	C9H10N2	001197-19-9	59



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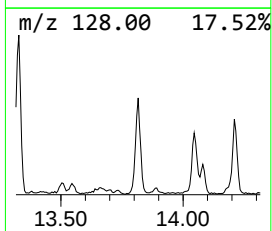
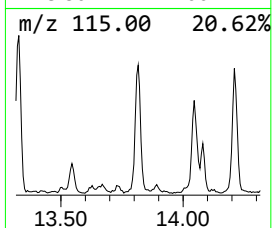
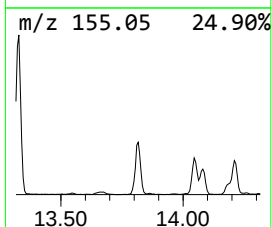
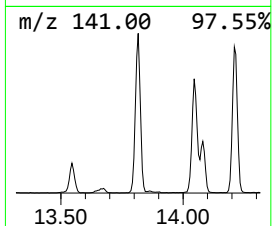
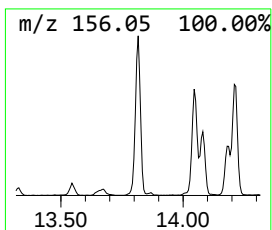
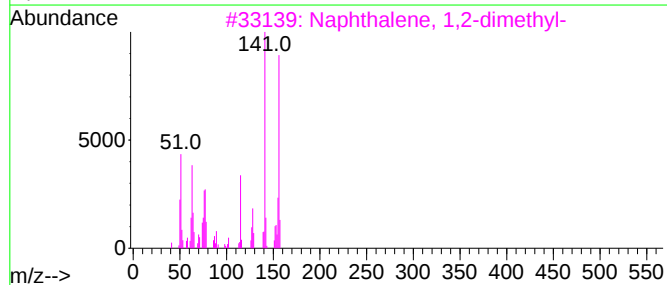
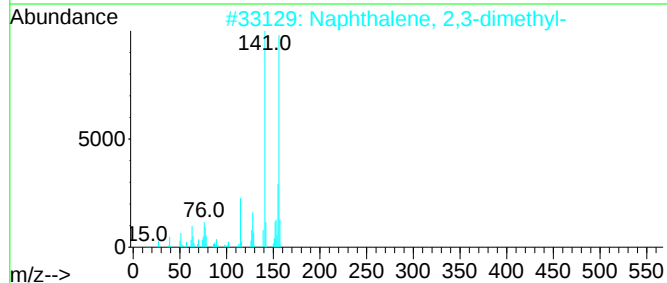
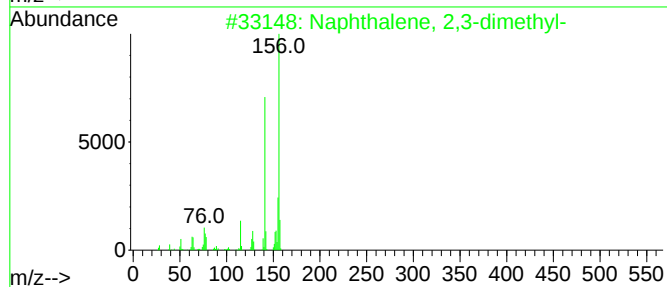
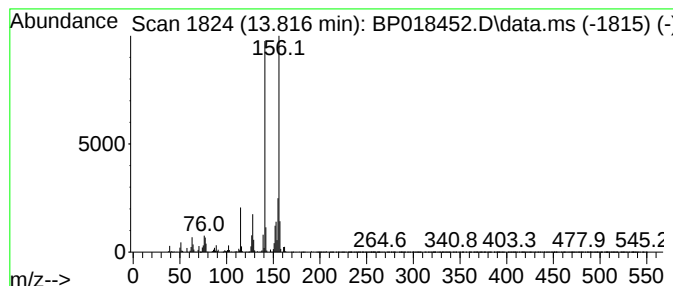
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Naphthalene, 2,3-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.816	4.60 ng/ul	729046	Acenaphthene-d10	14.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3			Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	97
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
5			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	96



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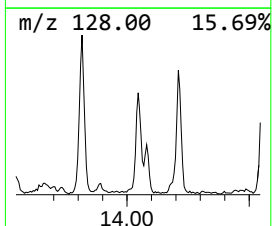
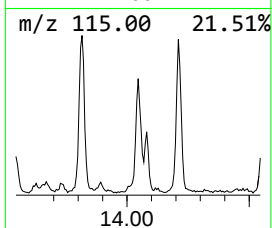
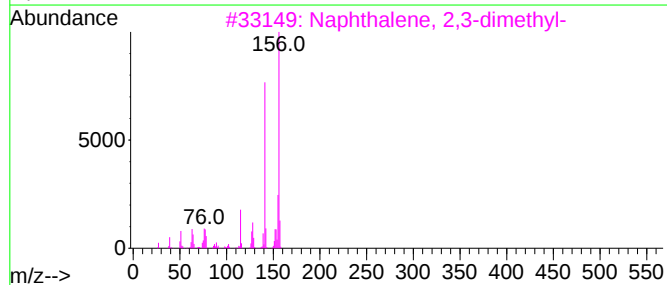
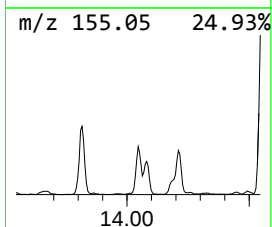
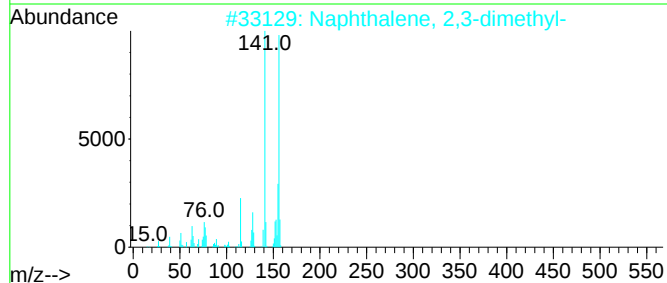
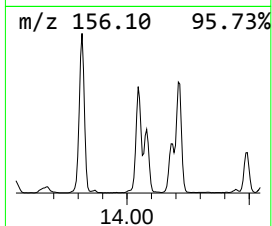
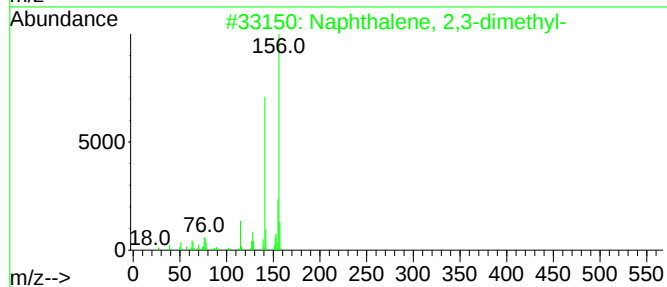
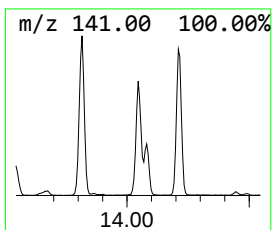
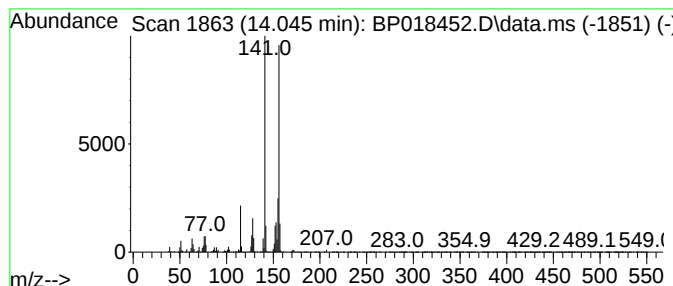
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 Naphthalene, 1,4-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.045	3.00 ng/ul	475809	Acenaphthene-d10	14.486

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
2			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
3			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
5			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112723\
Data File : BP018452.D
Acq On : 28 Nov 2023 22:04
Operator : MA/JU
Sample : 05332-16DL 5X
Misc :
ALS Vial : 61 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
YCG97DL

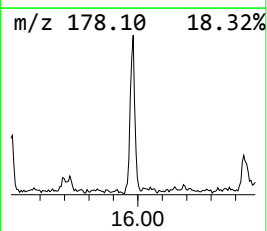
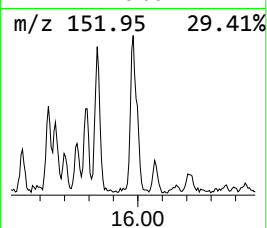
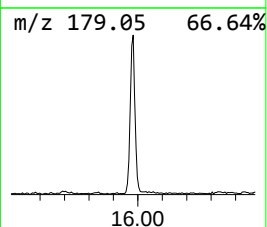
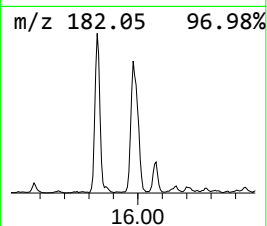
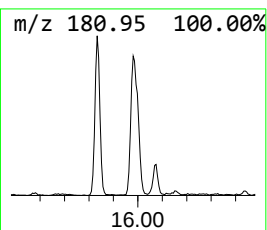
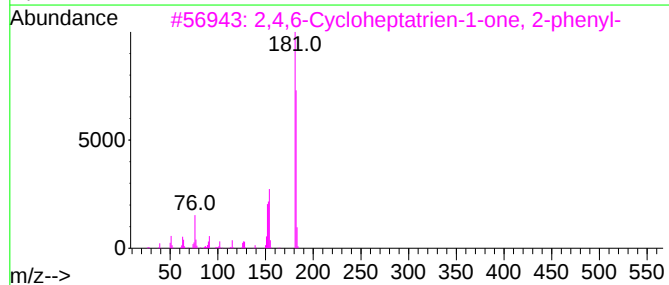
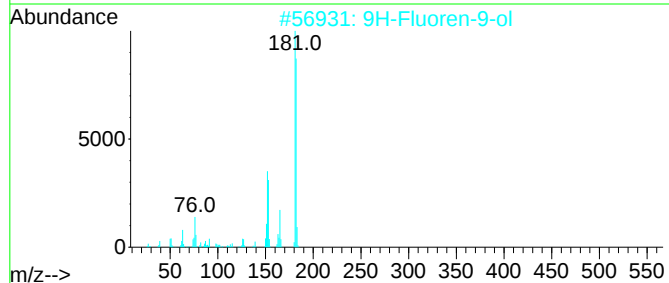
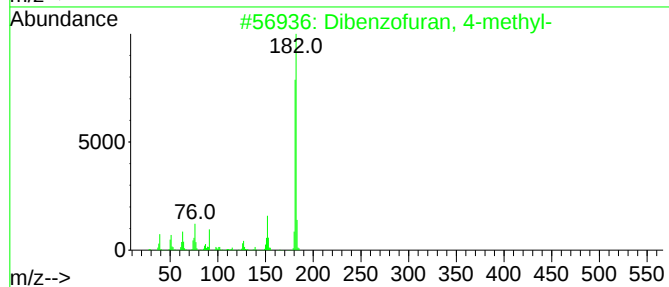
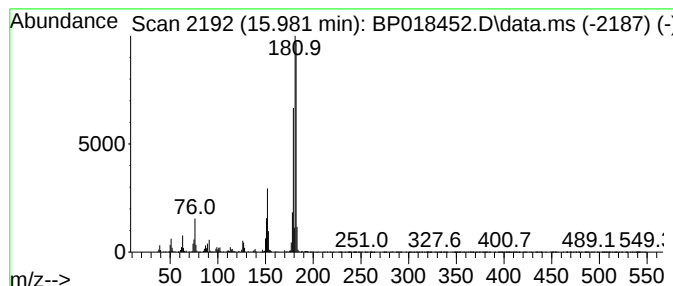
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Dibenzofuran, 4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.981	2.21 ng/ul	377759	Phenanthrene-d10	17.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzofuran, 4-methyl-	182	C13H10O	007320-53-8	72
2			9H-Fluoren-9-ol	182	C13H10O	001689-64-1	68
3			2,4,6-Cycloheptatrien-1-one, 2-p...	182	C13H10O	014562-09-5	62
4			2-Hydroxyfluorene	182	C13H10O	002443-58-5	53
5			3-(1-Naphthyl)acrylaldehyde	182	C13H10O	001504-73-0	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112723\
Data File : BP018452.D
Acq On : 28 Nov 2023 22:04
Operator : MA/JU
Sample : 05332-16DL 5X
Misc :
ALS Vial : 61 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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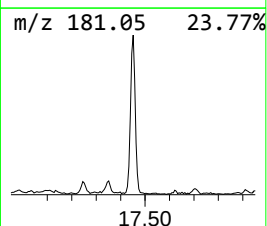
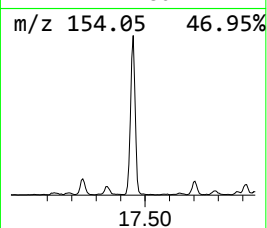
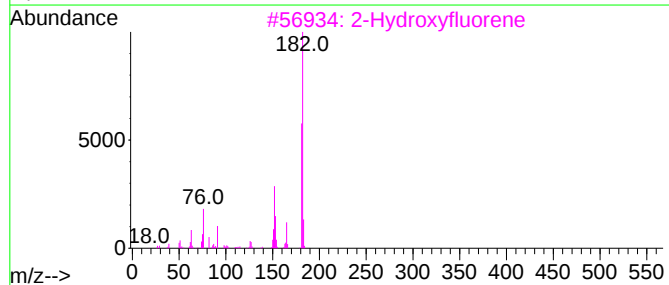
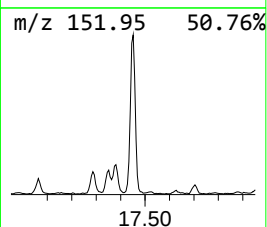
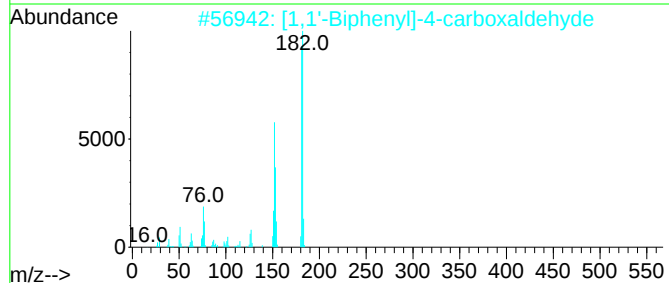
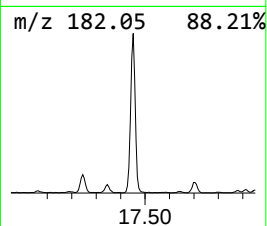
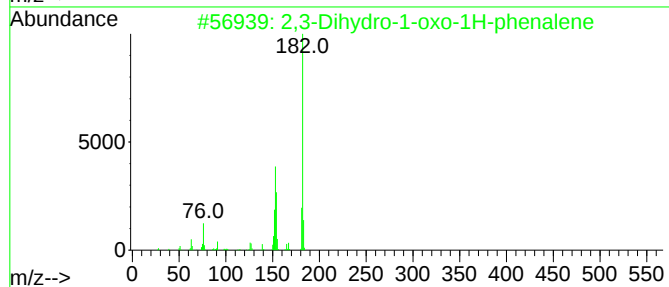
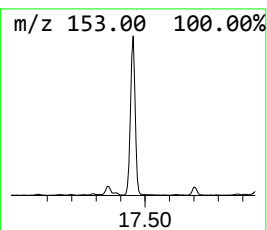
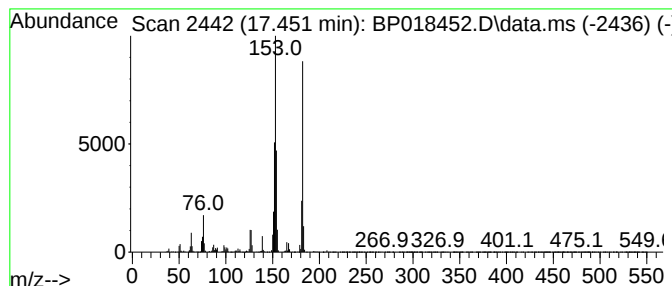
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 2,3-Dihydro-1-oxo-1H-phenalene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.451	5.82 ng/ul	995404	Phenanthrene-d10	17.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,3-Dihydro-1-oxo-1H-phenalene	182	C13H10O	000518-85-4	76
2			[1,1'-Biphenyl]-4-carboxaldehyde	182	C13H10O	003218-36-8	60
3			2-Hydroxyfluorene	182	C13H10O	002443-58-5	53
4			1,2,4-Cyclopentanetrione, 3-meth...	182	C9H10O4	057174-14-8	50
5			Anthracene, 1,2,3,4-tetrahydro-	182	C14H14	002141-42-6	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112723\
Data File : BP018452.D
Acq On : 28 Nov 2023 22:04
Operator : MA/JU
Sample : 05332-16DL 5X
Misc :
ALS Vial : 61 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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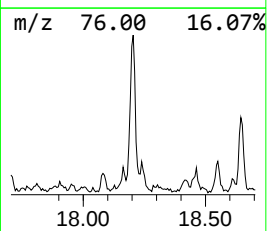
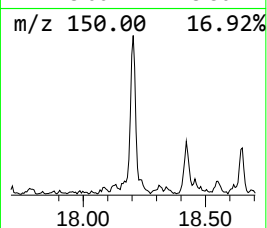
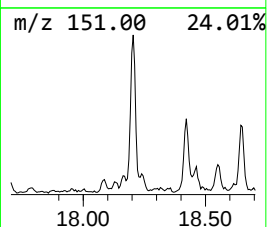
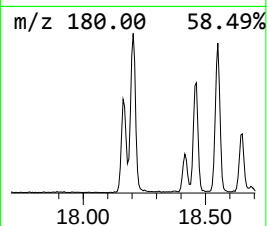
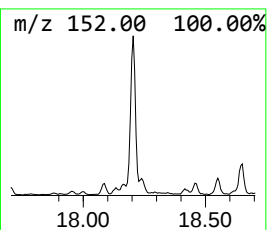
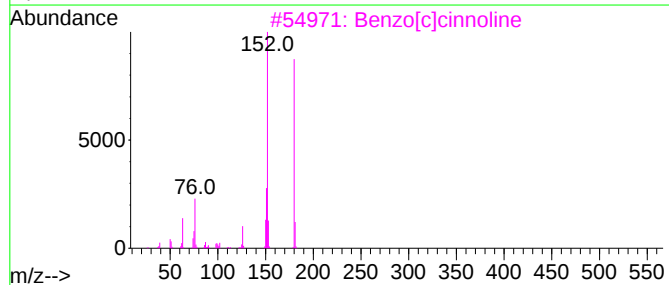
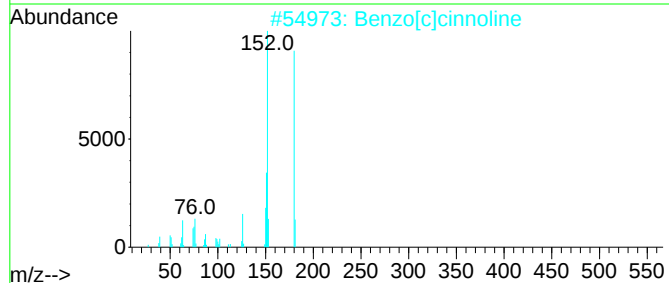
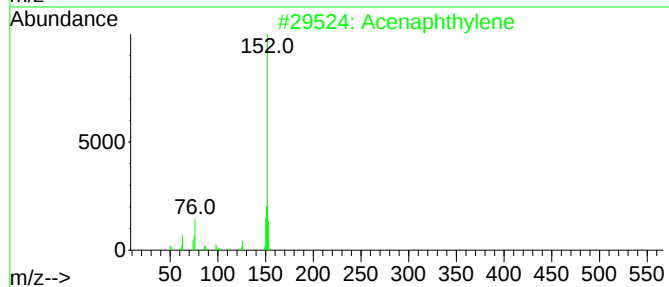
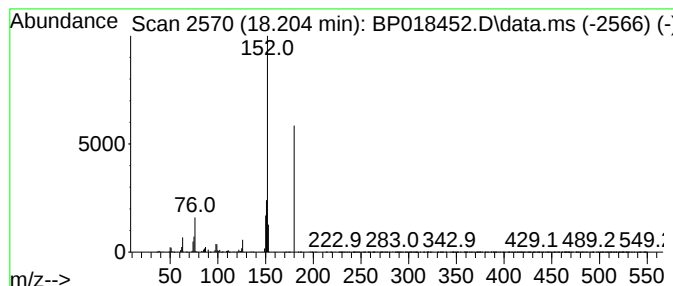
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzo[c]cinnoline Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.204	2.07 ng/ul	353513	Phenanthrene-d10	17.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acenaphthylene	152	C12H8	000208-96-8	95
2			Benzo[c]cinnoline	180	C12H8N2	000230-17-1	91
3			Benzo[c]cinnoline	180	C12H8N2	000230-17-1	91
4			Benzo[c]cinnoline	180	C12H8N2	000230-17-1	83
5			Biphenylene	152	C12H8	000259-79-0	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112723\
Data File : BP018452.D
Acq On : 28 Nov 2023 22:04
Operator : MA/JU
Sample : 05332-16DL 5X
Misc :
ALS Vial : 61 Sample Multiplier: 1

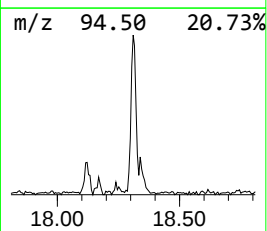
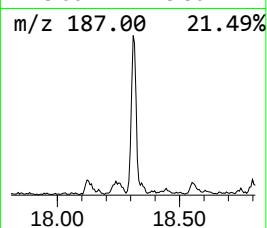
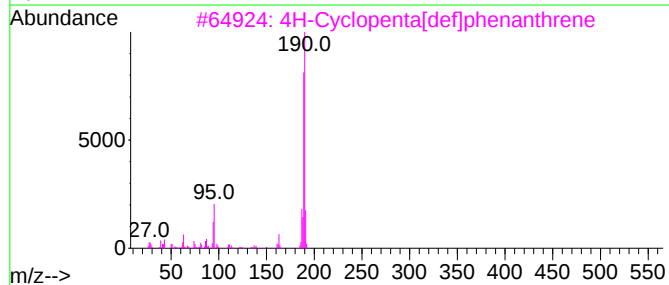
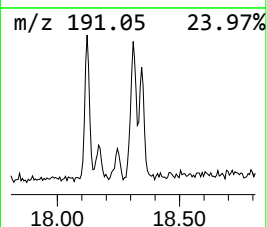
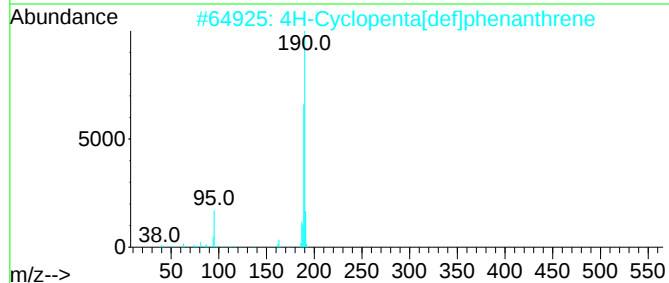
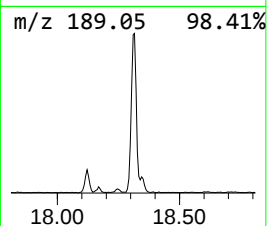
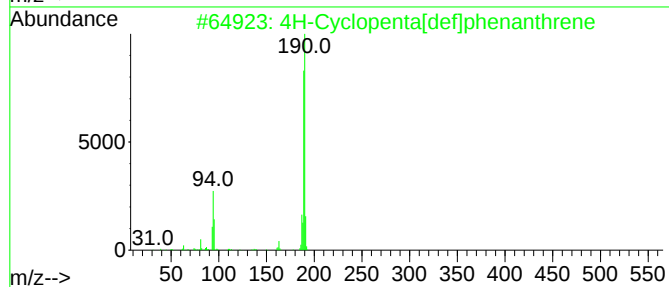
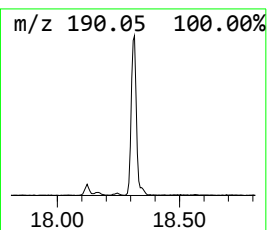
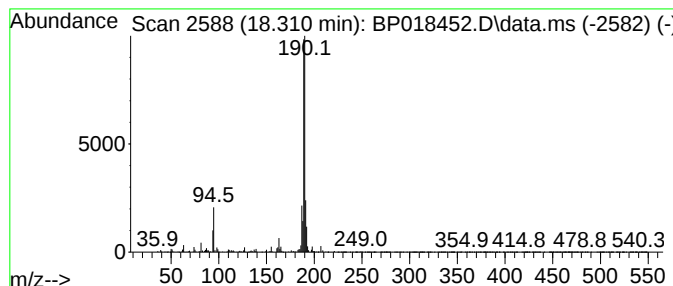
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 4H-Cyclopenta[def]phenanthrene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD			R.T.
18.310	2.48 ng/ul	424539	Phenanthrene-d10			17.245
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene		190	C15H10	000203-64-5	87
2	4H-Cyclopenta[def]phenanthrene		190	C15H10	000203-64-5	87
3	4H-Cyclopenta[def]phenanthrene		190	C15H10	000203-64-5	76
4	6H-Cyclobuta[jk]phenanthrene		190	C15H10	083469-43-6	72
5	Methyl diselenide		190	C2H6Se2	007101-31-7	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112723\
Data File : BP018452.D
Acq On : 28 Nov 2023 22:04
Operator : MA/JU
Sample : 05332-16DL 5X
Misc :
ALS Vial : 61 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112223.MA.M
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
Benzofuran, 4,7...	11.075	2.1	ng/ul	303282	2	10.651	2817200	20.0
Naphthalene, 2,...	13.816	4.6	ng/ul	729046	3	14.486	3169930	20.0
Naphthalene, 1,...	14.045	3.0	ng/ul	475809	3	14.486	3169930	20.0
Dibenzofuran, 4...	15.981	2.2	ng/ul	377759	4	17.245	3422320	20.0
2,3-Dihydro-1-o...	17.451	5.8	ng/ul	995404	4	17.245	3422320	20.0
Benzo[c]cinnoline	18.204	2.1	ng/ul	353513	4	17.245	3422320	20.0
4H-Cyclopenta[d...	18.310	2.5	ng/ul	424539	4	17.245	3422320	20.0