

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071824\
 Data File : BP021071.D
 Acq On : 19 Jul 2024 09:17
 Operator : MA/JU
 Sample : P3215-03DL 5X
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A4C00DL

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-BP071024.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP021071.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.017	3	5	15	rVB	128643	172663	3.25%	0.380%
2	3.228	39	41	50	rVB5	10390	17686	0.33%	0.039%
3	3.646	108	112	116	rVV	35727	46809	0.88%	0.103%
4	3.693	116	120	123	rVB5	5221	8473	0.16%	0.019%
5	3.740	123	128	135	rBV	169449	245120	4.61%	0.539%
6	3.817	138	141	146	rVB	24067	30793	0.58%	0.068%
7	3.952	160	164	171	rVB5	9629	19258	0.36%	0.042%
8	4.122	189	193	198	rVB6	6623	11051	0.21%	0.024%
9	4.499	252	257	263	rVB	40180	59683	1.12%	0.131%
10	4.652	281	283	286	rBV4	4760	5859	0.11%	0.013%
11	4.846	311	316	321	rVB2	12296	19026	0.36%	0.042%
12	4.958	331	335	345	rBV2	21568	42541	0.80%	0.094%
13	5.234	379	382	386	rBV4	4402	6283	0.12%	0.014%
14	6.905	660	666	678	rBV	74073	156766	2.95%	0.345%
15	7.028	681	687	703	rVB	87163	143488	2.70%	0.315%
16	7.240	714	723	729	rBV	100491	182722	3.44%	0.402%
17	7.681	791	798	811	rBV	579833	935159	17.58%	2.056%
18	8.428	920	925	930	rBV5	13965	30075	0.57%	0.066%
19	8.505	935	938	949	rVB9	7470	16197	0.30%	0.036%
20	8.846	987	996	1003	rBV	82896	180573	3.40%	0.397%
21	9.546	1109	1115	1128	rBV	84065	159170	2.99%	0.350%
22	9.828	1156	1163	1168	rBV10	3705	10637	0.20%	0.023%
23	10.116	1203	1212	1227	rBV	67870	197538	3.71%	0.434%
24	10.446	1260	1268	1274	rBV	709020	1357413	25.52%	2.984%
25	10.499	1274	1277	1287	rVB	79917	139628	2.63%	0.307%
26	10.622	1293	1298	1308	rBV3	19484	45633	0.86%	0.100%
27	11.199	1388	1396	1404	rBV2	19851	55366	1.04%	0.122%
28	12.040	1529	1539	1544	rVB2	3724	9980	0.19%	0.022%
29	12.116	1544	1552	1561	rBV	28695	55780	1.05%	0.123%
30	12.328	1581	1588	1595	rVB	16560	29917	0.56%	0.066%
31	12.787	1663	1666	1671	rVB6	5579	7972	0.15%	0.018%
32	13.134	1720	1725	1732	rBV3	8235	16085	0.30%	0.035%
33	13.328	1754	1758	1760	rBV5	5919	7898	0.15%	0.017%
34	13.481	1776	1784	1789	rBV8	8048	25594	0.48%	0.056%
35	13.628	1801	1809	1812	rBV3	15547	29935	0.56%	0.066%
36	13.675	1812	1817	1819	rVV5	11018	16272	0.31%	0.036%

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Integrator: RTE

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Peak Location: TOP

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
 Title : SVOA CALIBRATION

37	13.716	1819	1824	1835	rVB	273756	384257	7.23%	0.845%
38	13.869	1845	1850	1853	rBV5	6378	10565	0.20%	0.023%
39	13.963	1861	1866	1868	rBV6	5774	9155	0.17%	0.020%
40	14.004	1868	1873	1885	rVB	344624	488033	9.18%	1.073%
41	14.192	1899	1905	1908	rBV8	3263	6850	0.13%	0.015%
42	14.316	1914	1926	1932	rBV	1211112	1841721	34.63%	4.049%
43	14.375	1932	1936	1945	rVB	165707	252511	4.75%	0.555%
44	14.475	1949	1953	1954	rBV2	8125	9443	0.18%	0.021%
45	14.528	1954	1962	1972	rVB	529845	1138163	21.40%	2.502%
46	14.628	1972	1979	1988	rBV2	45070	111802	2.10%	0.246%
47	14.716	1988	1994	2002	rBV	105783	181587	3.41%	0.399%
48	14.934	2027	2031	2036	rVB6	6758	11782	0.22%	0.026%
49	14.992	2037	2041	2044	rBV6	3936	6135	0.12%	0.013%
50	15.116	2056	2062	2064	rBV7	5857	9682	0.18%	0.021%
51	15.163	2066	2070	2074	rVV7	4621	8136	0.15%	0.018%
52	15.322	2091	2097	2102	rBV	443785	625339	11.76%	1.375%
53	15.381	2102	2107	2113	rVB	207115	313061	5.89%	0.688%
54	15.487	2120	2125	2132	rBV5	43608	99919	1.88%	0.220%
55	15.557	2133	2137	2140	rVV2	18341	30883	0.58%	0.068%
56	15.592	2141	2143	2148	rVV6	10155	16971	0.32%	0.037%
57	15.639	2148	2151	2153	rVV3	11890	15483	0.29%	0.034%
58	15.675	2153	2157	2164	rVB	32000	46269	0.87%	0.102%
59	15.839	2180	2185	2192	rBV	31365	61599	1.16%	0.135%
60	16.086	2222	2227	2237	rVB8	18670	43232	0.81%	0.095%
61	16.275	2255	2259	2263	rBV7	4918	8844	0.17%	0.019%
62	16.404	2274	2281	2288	rBV6	25581	66178	1.24%	0.145%
63	16.463	2288	2291	2295	rVB5	14245	19508	0.37%	0.043%
64	16.639	2314	2321	2326	rBV8	16422	41643	0.78%	0.092%
65	16.686	2326	2329	2332	rVV5	8083	13480	0.25%	0.030%
66	16.786	2340	2346	2350	rBV	70689	114934	2.16%	0.253%
67	16.839	2352	2355	2358	rBV2	16761	25770	0.48%	0.057%
68	16.939	2367	2372	2379	rVB	148940	214485	4.03%	0.472%
69	17.128	2396	2404	2407	rBV	1306900	2335734	43.92%	5.135%
70	17.175	2407	2412	2417	rVV	2538026	3978549	74.81%	8.746%
71	17.228	2417	2421	2424	rVV2	520279	766374	14.41%	1.685%
72	17.257	2424	2426	2433	rVB	579471	749746	14.10%	1.648%
73	17.551	2468	2476	2483	rBV2	275444	508066	9.55%	1.117%
74	17.657	2490	2494	2497	rBV	23363	31847	0.60%	0.070%
75	17.857	2526	2528	2536	rVB3	31063	48535	0.91%	0.107%
76	18.022	2552	2556	2559	rBV	153374	213844	4.02%	0.470%

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77	18.075	2559	2565	2572	rVB2	215517	452230	8.50%	0.994%
78	18.163	2577	2580	2585	rVB	56410	81790	1.54%	0.180%
79	18.233	2586	2592	2596	rBV2	357103	589271	11.08%	1.295%
80	18.269	2596	2598	2604	rVB	102540	113795	2.14%	0.250%
81	18.345	2609	2611	2617	rBV6	23817	35782	0.67%	0.079%
82	18.551	2641	2646	2651	rBV	139198	213705	4.02%	0.470%
83	18.610	2651	2656	2662	rBV	167456	262696	4.94%	0.577%
84	18.669	2664	2666	2669	rBV3	22850	28394	0.53%	0.062%
85	19.033	2725	2728	2734	rBV4	62884	98023	1.84%	0.215%
86	19.139	2743	2746	2750	rVB	82810	99677	1.87%	0.219%
87	19.257	2760	2766	2774	rVB	3786112	5318382	100.00%	11.692%
88	19.598	2819	2824	2827	rBV3	959427	1557525	29.29%	3.424%
89	19.633	2827	2830	2839	rVV	2050344	2707806	50.91%	5.953%
90	19.716	2840	2844	2848	rVB2	90723	141654	2.66%	0.311%
91	19.833	2860	2864	2868	rBV	137111	178738	3.36%	0.393%
92	19.898	2872	2875	2876	rBV	83125	79268	1.49%	0.174%
93	20.157	2916	2919	2928	rVB	268221	399510	7.51%	0.878%
94	20.257	2933	2936	2940	rBV2	188202	257802	4.85%	0.567%
95	21.133	3081	3085	3089	rBV2	300520	556161	10.46%	1.223%
96	21.569	3151	3159	3163	rBV2	2197861	5299748	99.65%	11.651%
97	21.610	3163	3166	3179	rVB	1365742	2220031	41.74%	4.880%
98	23.833	3537	3544	3550	rBV2	813938	2315301	43.53%	5.090%
99	24.863	3713	3719	3730	rVB	1187209	3126714	58.79%	6.874%

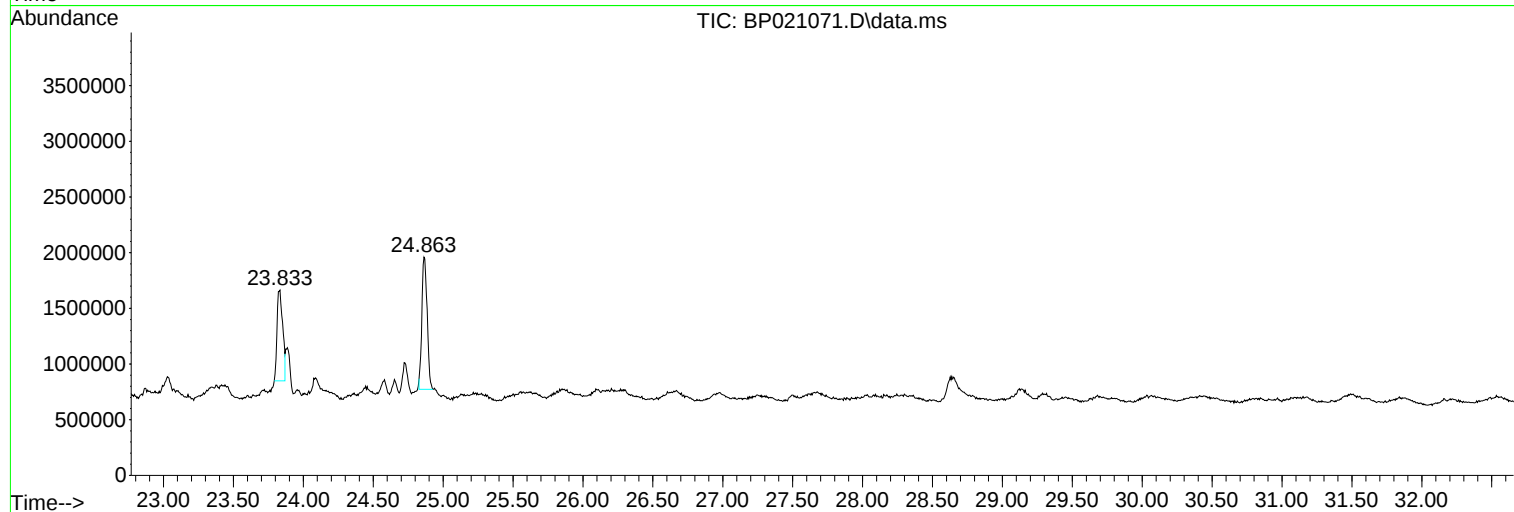
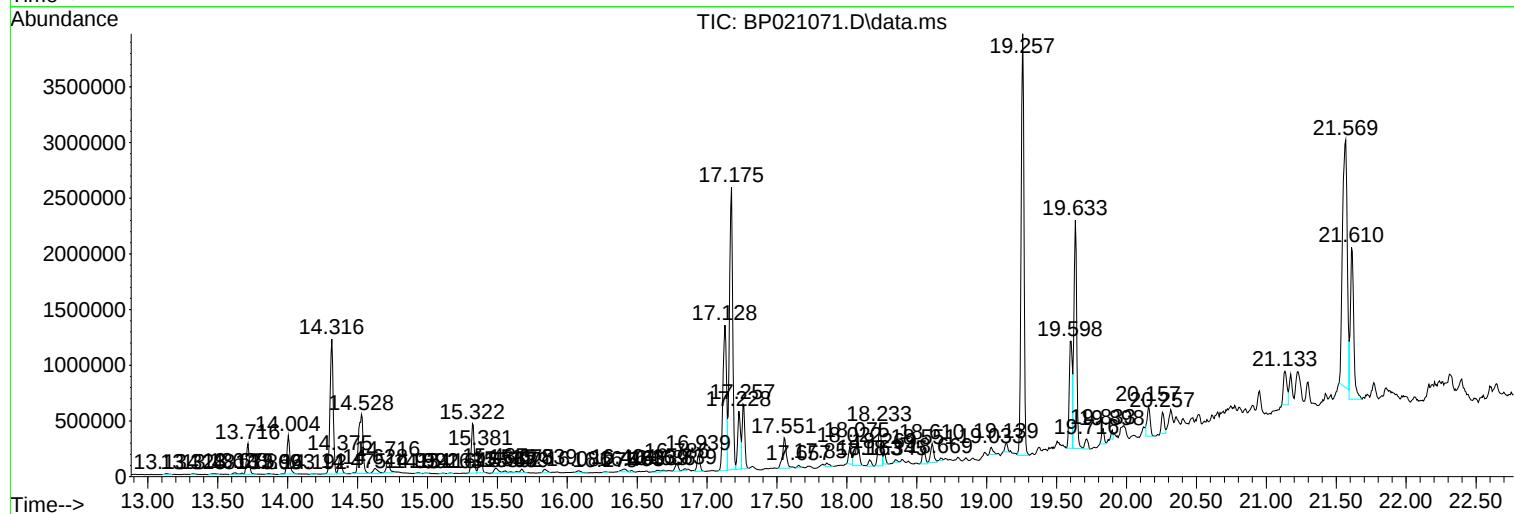
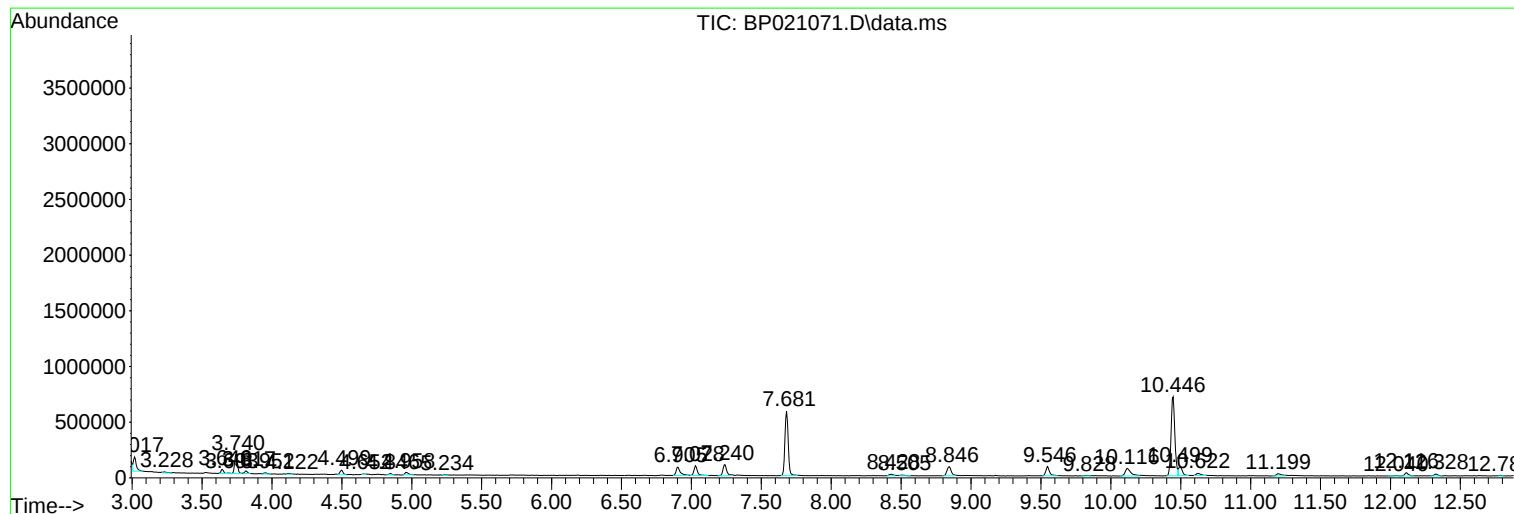
Sum of corrected areas: 45489161

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ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4C00DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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



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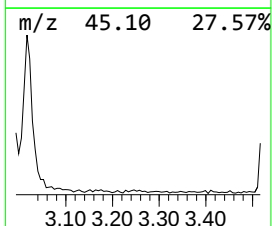
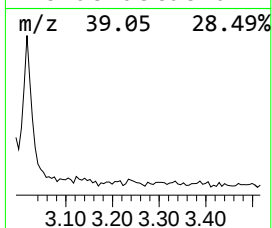
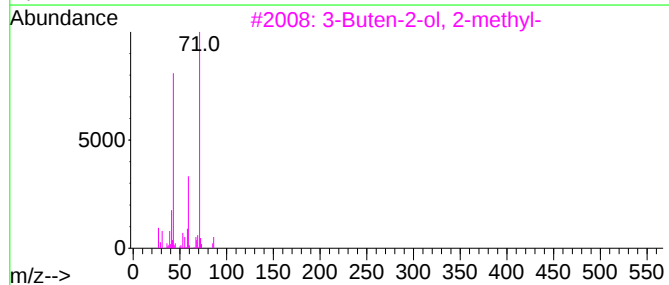
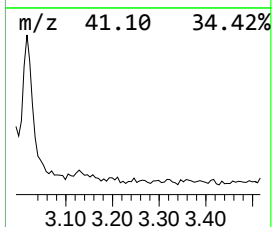
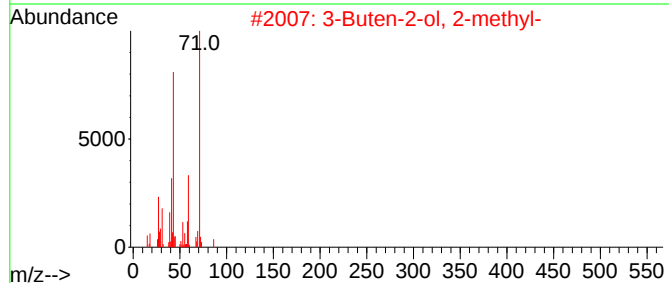
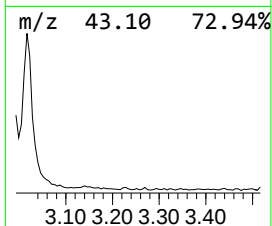
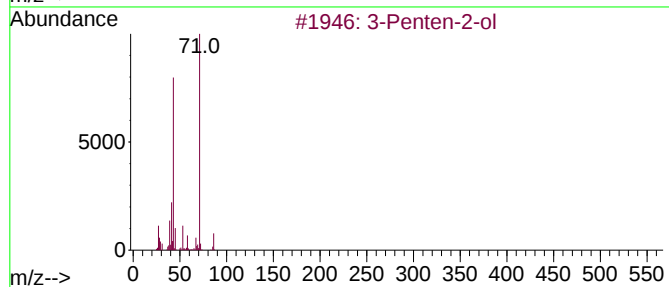
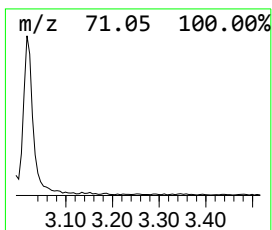
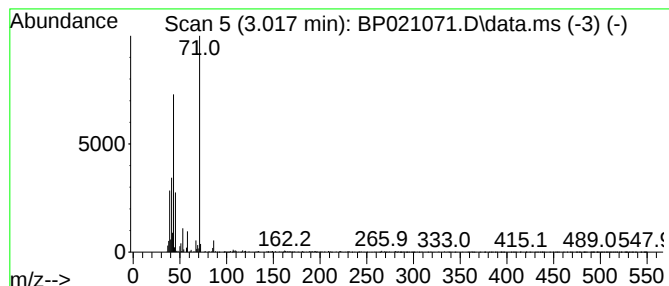
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 3-Penten-2-ol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.017	3.69 ng/ul	172663	1,4-Dichlorobenzene-d4	7.681

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Penten-2-ol	86	C5H10O	001569-50-2	72
2			3-Buten-2-ol, 2-methyl-	86	C5H10O	000115-18-4	64
3			3-Buten-2-ol, 2-methyl-	86	C5H10O	000115-18-4	59
4			2-Methylpropionic acid, 4-nitrop...	209	C10H11NO4	004195-16-8	47
5			Furan-2-carbonyl chloride, tetra...	134	C5H7ClO2	052449-98-6	47



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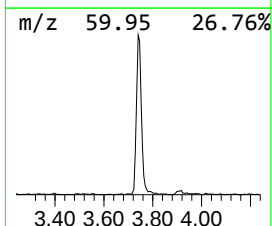
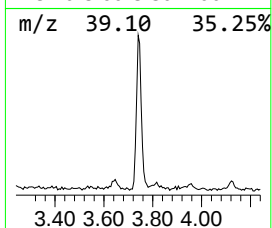
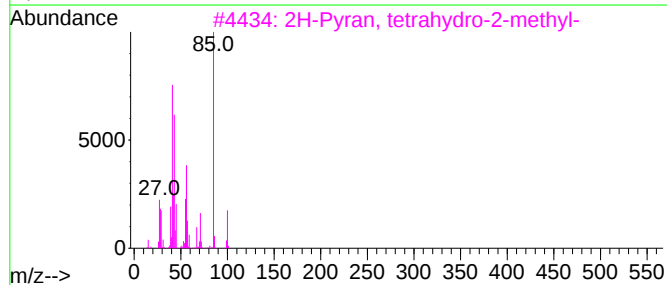
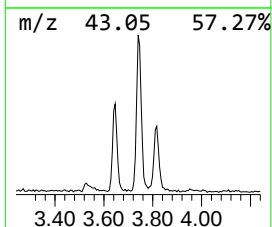
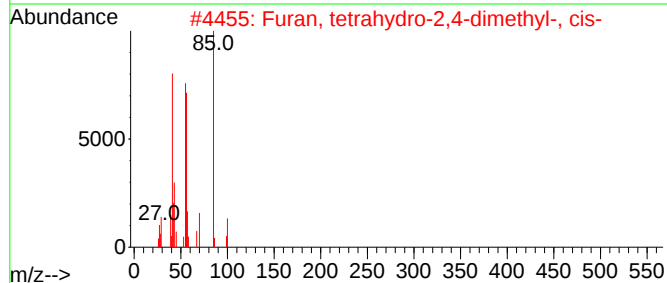
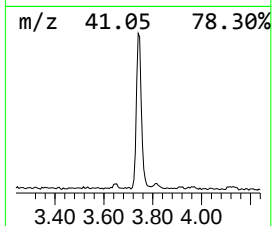
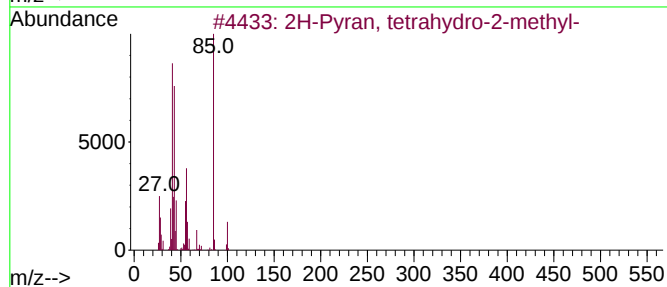
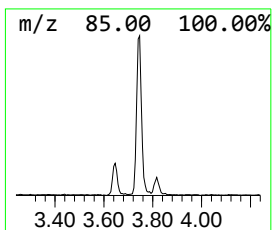
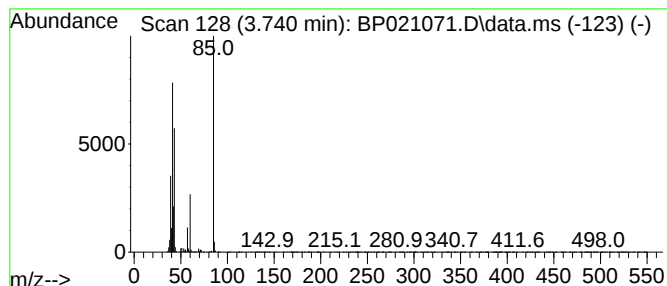
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2H-Pyran, tetrahydro-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.740	5.24 ng/ul	245120	1,4-Dichlorobenzene-d4	7.681

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	64
2			Furan, tetrahydro-2,4-dimethyl-,...	100	C6H12O	039168-01-9	38
3			2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	38
4			Hexane, 2-bromo-	164	C6H13Br	003377-86-4	36
5			Oxirane, (2-methylpropyl)-	100	C6H12O	023850-78-4	36



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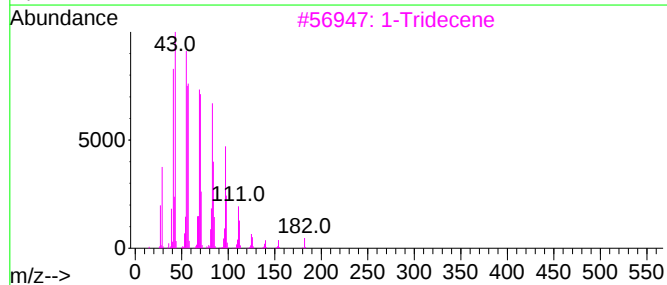
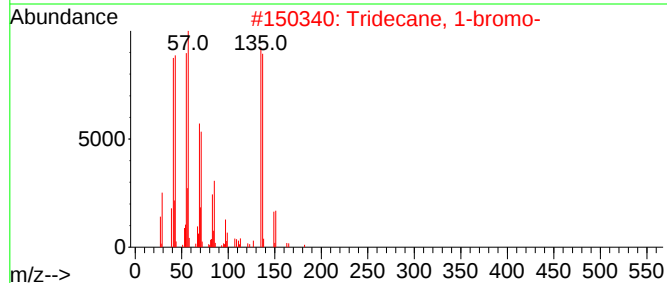
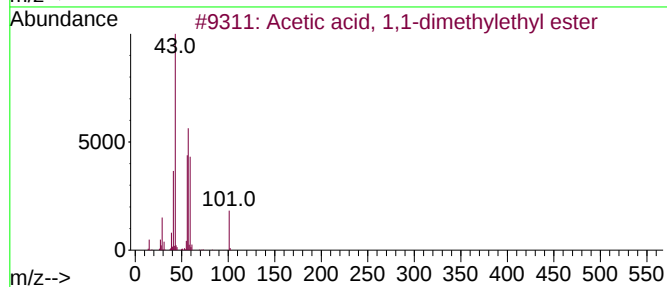
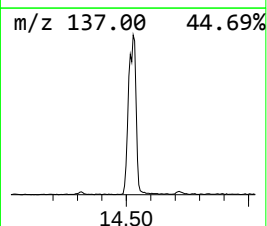
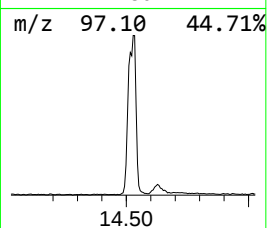
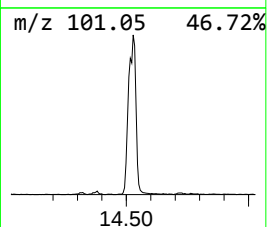
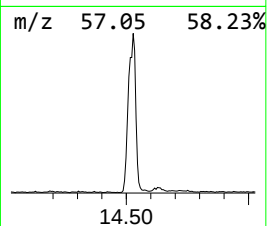
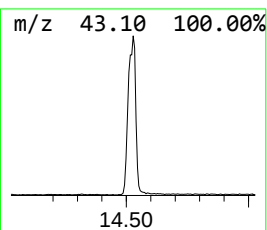
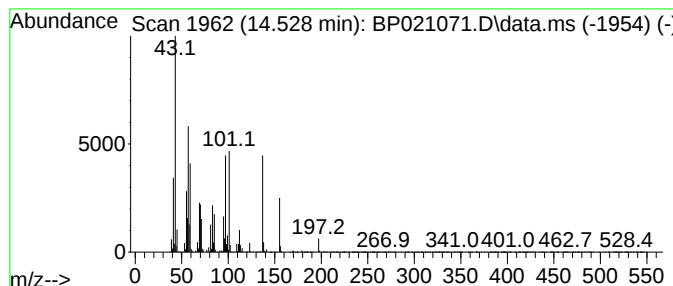
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ClientSampleId :
A4C00DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.528	12.36 ng/ul	1138160	Acenaphthene-d10		14.316	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Acetic acid, 1,1-dimethylethyl e...		116	C6H12O2	000540-88-5	10
2	Tridecane, 1-bromo-		262	C13H27Br	000765-09-3	10
3	1-Tridecene		182	C13H26	002437-56-1	10
4	11-Tricosene		322	C23H46	052078-56-5	10
5	Tetradecane, 1-bromo-		276	C14H29Br	000112-71-0	10



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071824\
Data File : BP021071.D
Acq On : 19 Jul 2024 09:17
Operator : MA/JU
Sample : P3215-03DL 5X
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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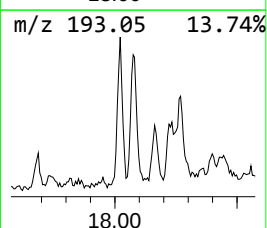
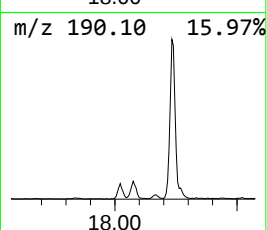
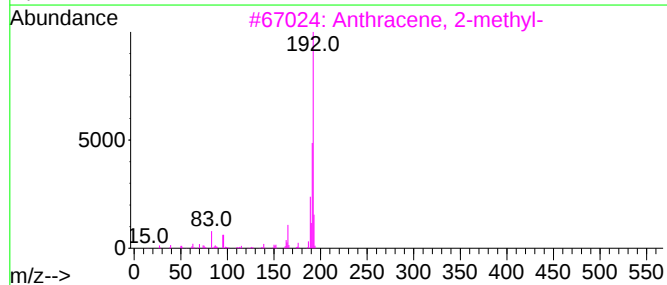
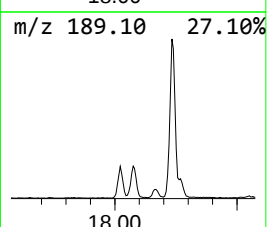
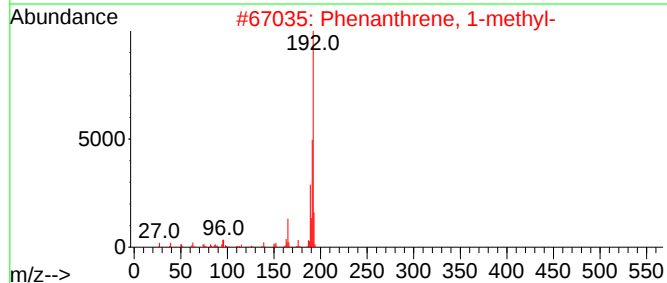
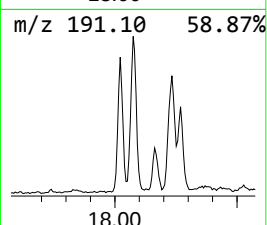
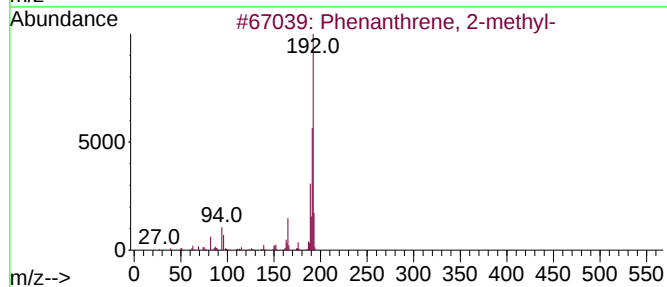
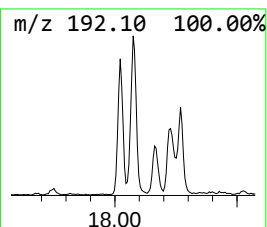
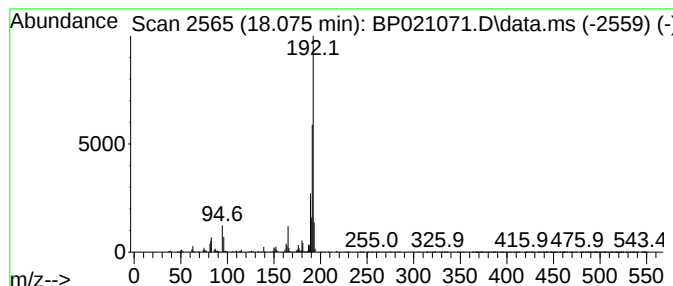
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Phenanthrene, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.075	3.87 ng/ul	452230	Phenanthrene-d10	17.128

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	98
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	96
3			Anthracene, 2-methyl-	192	C15H12	000613-12-7	96
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
5			1H-Cyclopropa[1]phenanthrene,1a,...	192	C15H12	000949-41-7	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071824\
Data File : BP021071.D
Acq On : 19 Jul 2024 09:17
Operator : MA/JU
Sample : P3215-03DL 5X
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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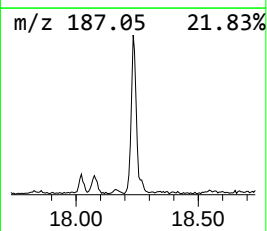
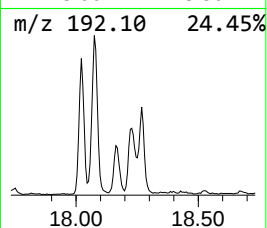
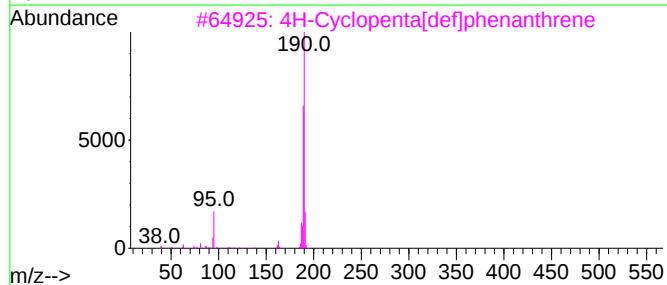
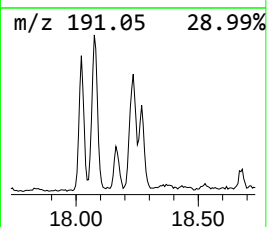
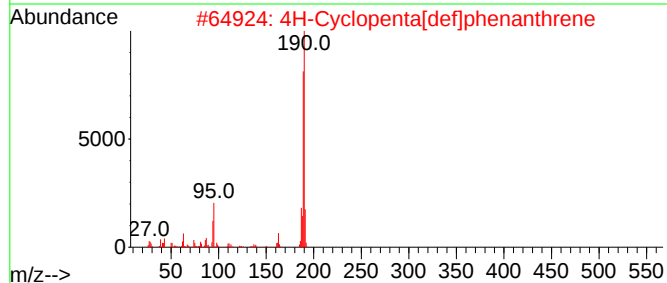
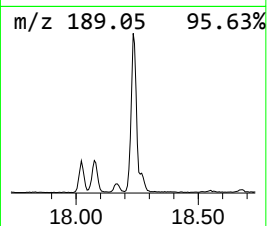
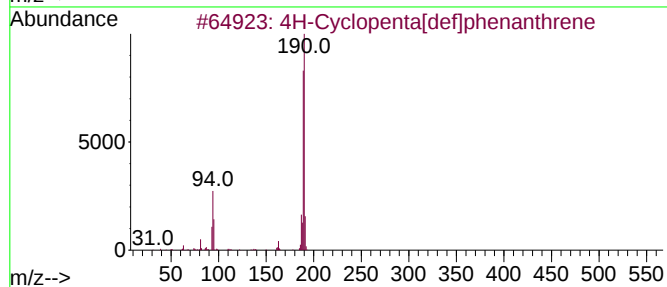
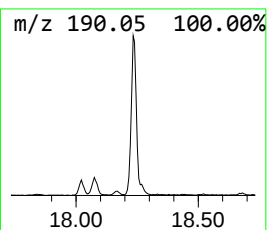
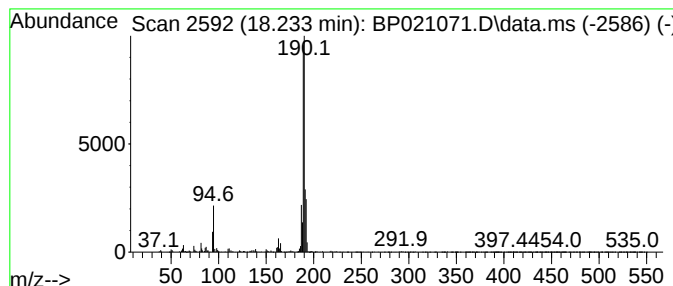
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.233	5.05 ng/ul	589271	Phenanthrene-d10	17.128

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	81
2			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	68
3			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	64
4			6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	50
5			Methyl diselenide	190	C2H6Se2	007101-31-7	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071824\
Data File : BP021071.D
Acq On : 19 Jul 2024 09:17
Operator : MA/JU
Sample : P3215-03DL 5X
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4C00DL

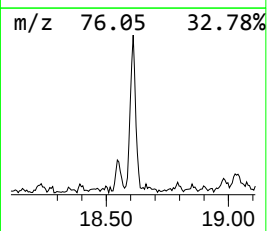
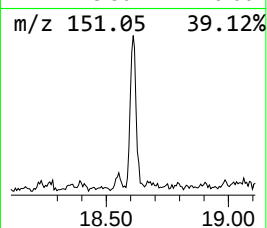
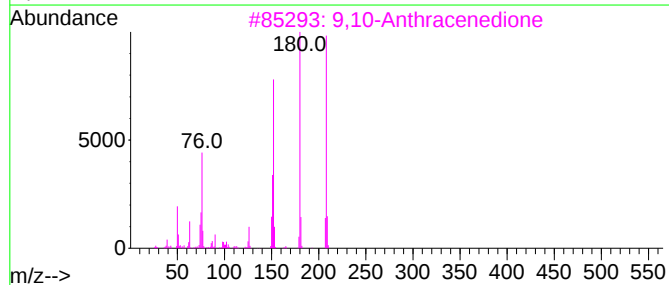
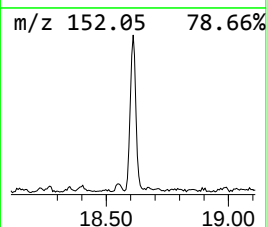
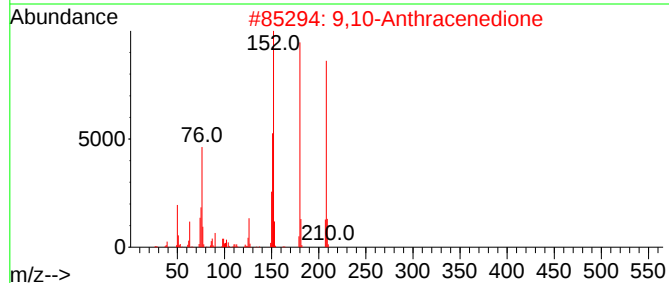
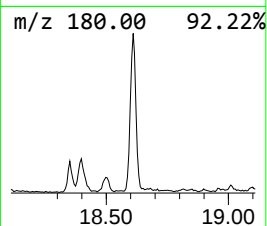
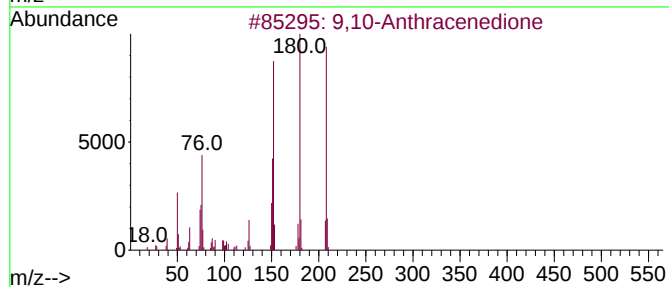
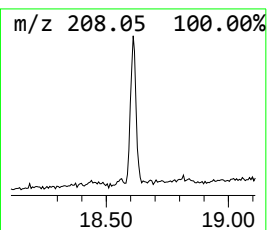
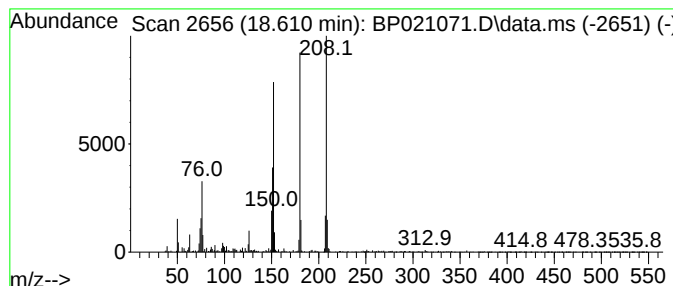
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 9,10-Anthracenedione Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.610	2.25 ng/ul	262696	Phenanthrene-d10	17.128

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione			208	C14H8O2	000084-65-1	98
2	9,10-Anthracenedione			208	C14H8O2	000084-65-1	98
3	9,10-Anthracenedione			208	C14H8O2	000084-65-1	95
4	9,10-Anthracenedione			208	C14H8O2	000084-65-1	94
5	Benzo[c]cinnoline			180	C12H8N2	000230-17-1	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071824\
Data File : BP021071.D
Acq On : 19 Jul 2024 09:17
Operator : MA/JU
Sample : P3215-03DL 5X
Misc :
ALS Vial : 30 Sample Multiplier: 1

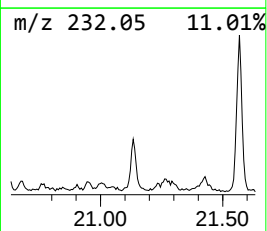
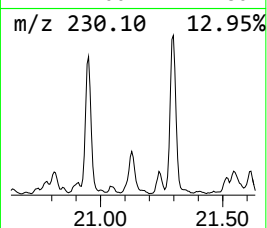
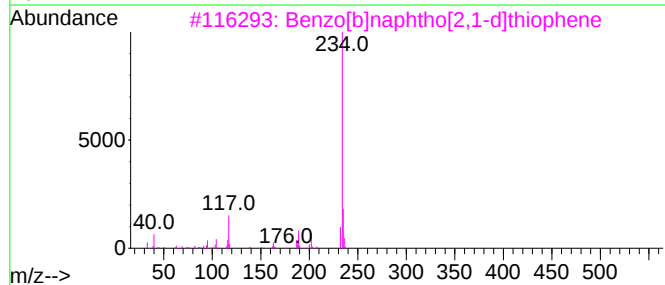
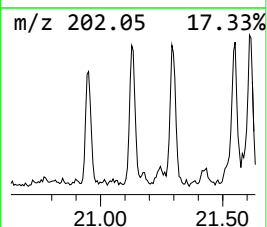
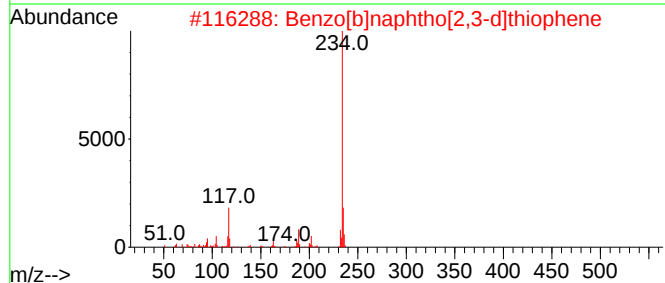
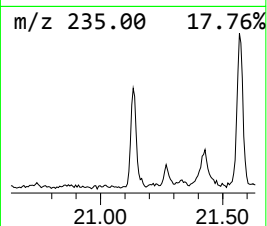
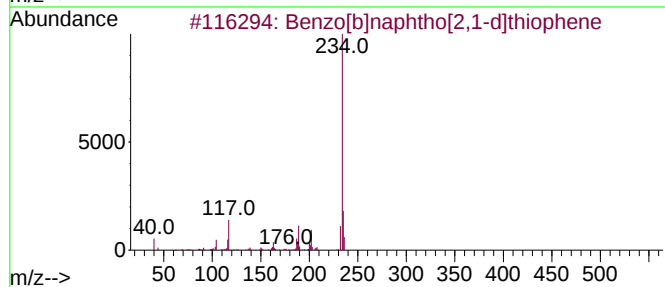
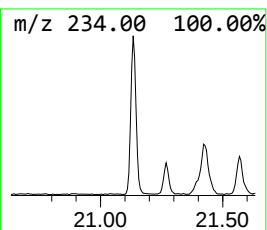
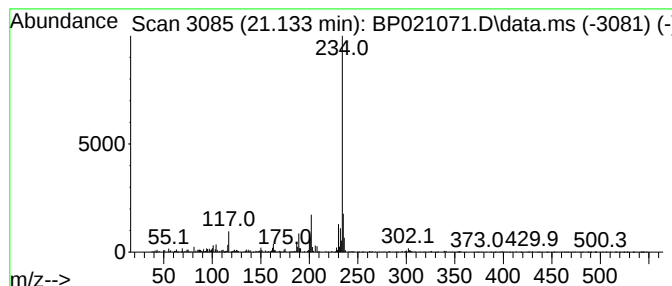
Instrument :
BNA_P
ClientSampleId :
A4C00DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzo[b]naphtho[2,1-d]thiop... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.133	2.10 ng/ul	556161	Chrysene-d12	21.569	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	94
2	Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	91
3	Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	90
4	Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	87
5	Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071824\
Data File : BP021071.D
Acq On : 19 Jul 2024 09:17
Operator : MA/JU
Sample : P3215-03DL 5X
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.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
3-Penten-2-ol	3.017	3.7	ng/u1	172663	1	7.681	935159	20.0
2H-Pyran, tetra...	3.740	5.2	ng/u1	245120	1	7.681	935159	20.0
unknown-01	14.528	12.4	ng/u1	1138160	3	14.316	1841720	20.0
Phenanthrene, 2...	18.075	3.9	ng/u1	452230	4	17.128	2335730	20.0
4H-Cyclopenta[d...	18.233	5.0	ng/u1	589271	4	17.128	2335730	20.0
9,10-Anthracene...	18.610	2.3	ng/u1	262696	4	17.128	2335730	20.0
Benzo[b]naphtho...	21.133	2.1	ng/u1	556161	5	21.569	5299750	20.0