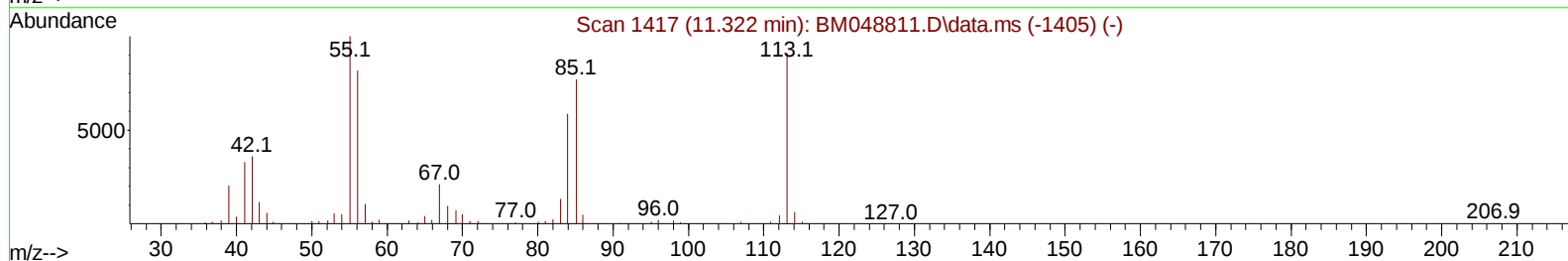
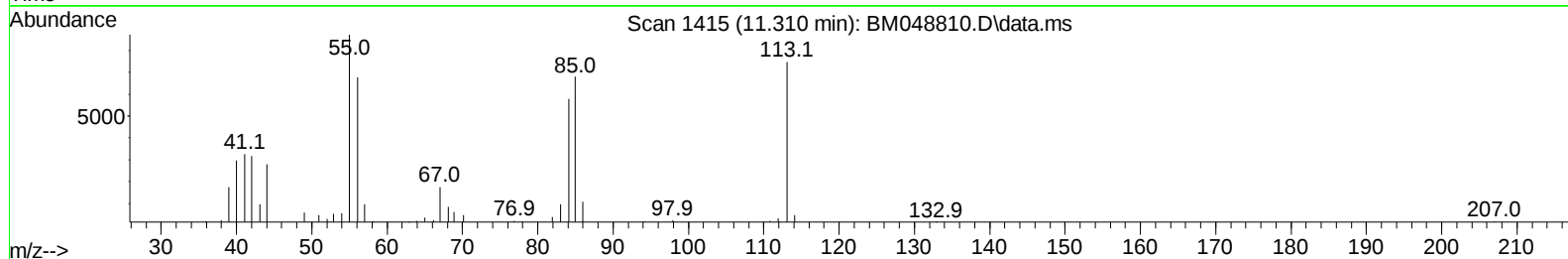
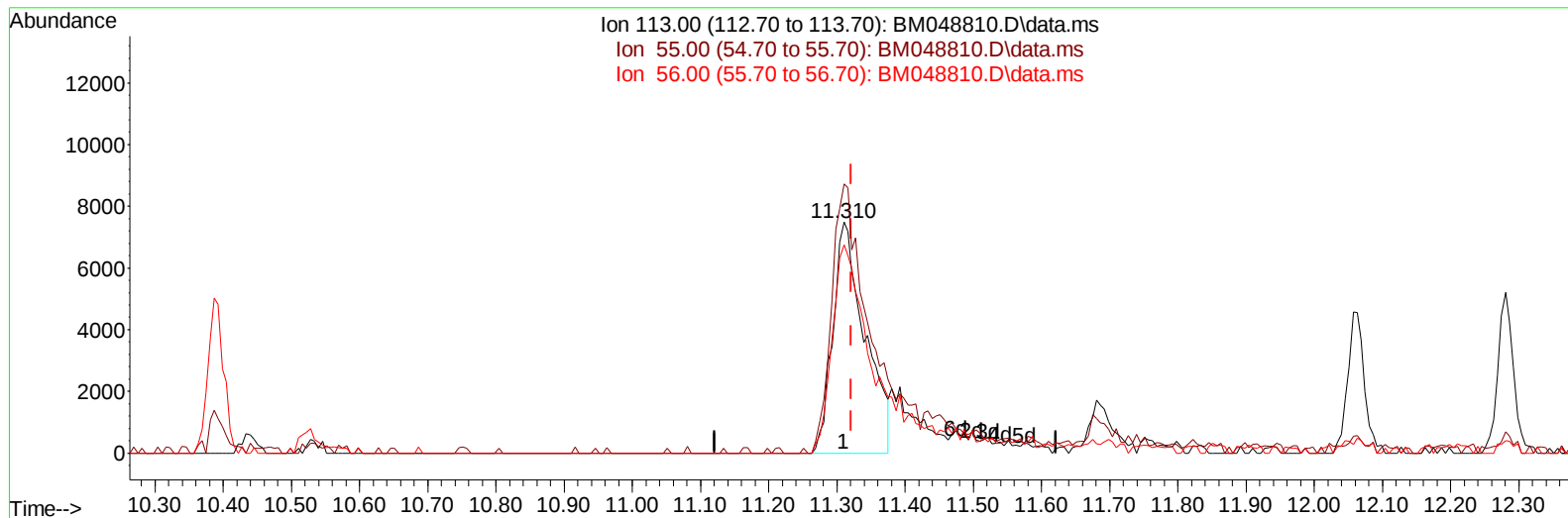


Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112224\
Data File : BM048810.D
Acq On : 22 Nov 2024 10:58
Operator : RC/JU
Sample : SST01061
Misc :
ALS Vial : 3 Sample Multiplier: 1

Quant Time: Nov 22 14:29:06 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112224.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Nov 22 14:18:42 2024
Response via : Initial Calibration



TIC: BM048810.D\data.ms

(34) Caprolactam

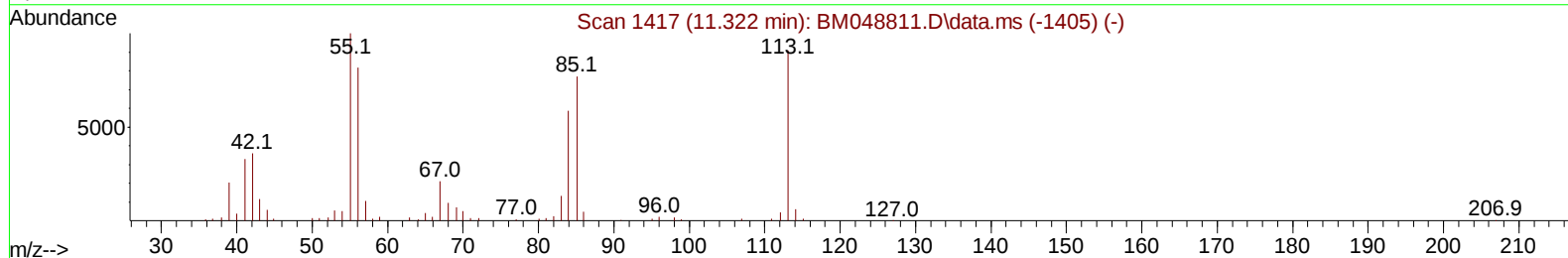
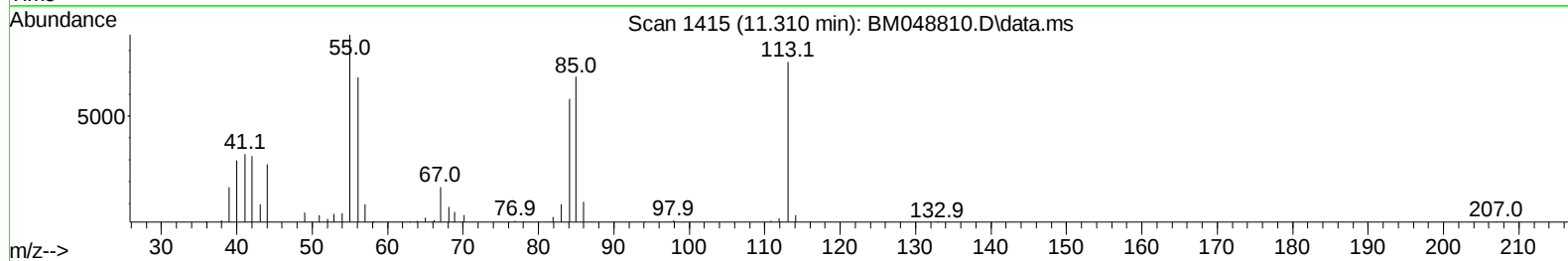
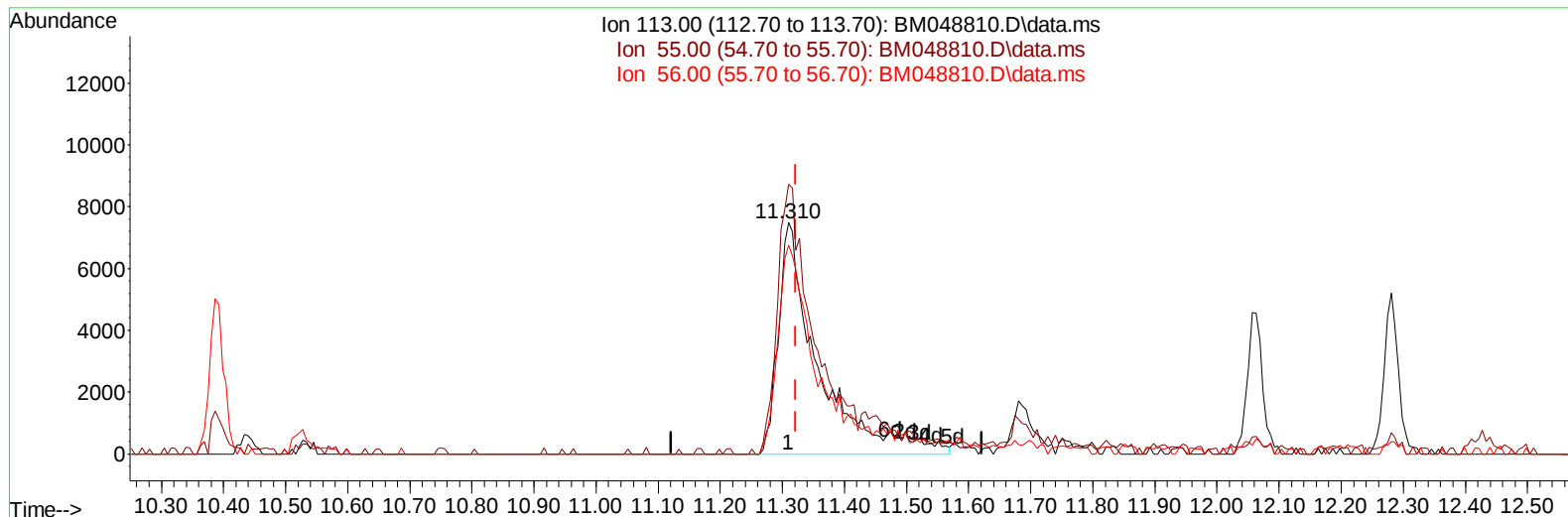
11.310min (-0.012) 6.29 ng/ul

response 24530

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	109.90	116.51
56.00	89.90	90.34
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112224\
Data File : BM048810.D
Acq On : 22 Nov 2024 10:58
Operator : RC/JU
Sample : SSTD01061
Misc :
ALS Vial : 3 Sample Multiplier: 1

Quant Time: Nov 22 14:29:06 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112224.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Nov 22 14:18:42 2024
Response via : Initial Calibration



TIC: BM048810.D\data.ms

(34) Caprolactam

11.310min (-0.012) 8.50 ng/ul m

response 33133

Ion	Exp%	Act%
113.00	100.00	100.00
55.00	109.90	116.51
56.00	89.90	90.34
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112224\
 Data File : BM048810.D
 Acq On : 22 Nov 2024 10:58
 Operator : RC/JU
 Sample : SSTD01061
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Nov 22 14:29:51 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 22 14:18:42 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.622	152	227092	20.000	ng/ul	0.00
20) Naphthalene-d8	10.392	136	979663	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.257	164	566716	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.004	188	1085166	20.000	ng/ul	0.00
79) Chrysene-d12	21.245	240	934907	20.000	ng/ul	0.00
88) Perylene-d12	24.103	264	993785	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.187	96	24289	4.128	ng/uL	0.00
4) Pyridine-d5	3.575	84	159555	9.468	ng/ul	0.00
7) Phenol-d5	6.793	99	154509	8.540	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.963	67	120308	9.726	ng/ul	0.00
11) 2-Chlorophenol-d4	7.157	132	123660	8.904	ng/ul	0.00
15) 4-Methylphenol-d8	8.322	113	116844	8.683	ng/ul	0.00
21) Nitrobenzene-d5	8.763	128	58226	8.871	ng/ul	0.00
24) 2-Nitrophenol-d4	9.481	143	28428	5.990	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.016	165	98378	8.150	ng/ul	0.00
31) 4-Chloroaniline-d4	10.528	131	183282	9.706	ng/ul	0.00
46) Dimethylphthalate-d6	13.680	166	345102	9.503	ng/ul	0.00
49) Acenaphthylene-d8	13.945	160	431313	9.549	ng/ul	0.00
54) 4-Nitrophenol-d4	14.457	143	30947	5.859	ng/ul	0.00
60) Fluorene-d10	15.257	176	299919	9.669	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.368	200	12097	7.218	ng/ul	0.00
73) Anthracene-d10	17.104	188	449566	9.631	ng/ul	0.00
81) Pyrene-d10	19.404	212	459910	9.044	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.909	264	407265	8.873	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.222	88	26619	4.233	ng/uL	97
5) Pyridine	3.593	79	176535	10.166	ng/ul	96
6) Benzaldehyde	6.769	77	127350	12.829	ng/ul	98
8) Phenol	6.822	94	171582	8.908	ng/ul	98
10) Bis(2-Chloroethyl)ether	7.057	93	153593	9.439	ng/ul	98
12) 2-Chlorophenol	7.187	128	137501	9.409	ng/ul	99
13) 2-Methylphenol	8.063	108	123994	8.841	ng/ul	98
14) 2,2'-oxybis(1-Chloropr...	8.145	45	104044	9.609	ng/ul	100
16) Acetophenone	8.434	105	229224	9.829	ng/ul	99
17) N-Nitroso-di-n-propyla...	8.422	70	119209	9.843	ng/ul	98
18) 4-Methylphenol	8.387	108	128741	8.612	ng/ul	97
19) Hexachloroethane	8.692	117	62873	9.417	ng/ul	95
22) Nitrobenzene	8.804	77	165400	9.337	ng/ul	97
23) Isophorone	9.328	82	309918	9.277	ng/ul	99
25) 2-Nitrophenol	9.510	139	39084	6.611	ng/ul	99
26) 2,4-Dimethylphenol	9.581	107	143860	9.529	ng/ul	99
27) Bis(2-Chloroethoxy)met...	9.822	93	190184	9.547	ng/ul	97
29) 2,4-Dichlorophenol	10.039	162	107932	8.910	ng/ul	94
30) Naphthalene	10.439	128	485955	9.930	ng/ul	99
32) 4-Chloroaniline	10.551	127	180766	9.883	ng/ul	96
33) Hexachlorobutadiene	10.739	225	78409	9.170	ng/ul	92
34) Caprolactam	11.310	113	33133m	8.500	ng/ul	
35) 4-Chloro-3-methylphenol	11.680	107	113027	8.852	ng/ul	96

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112224\
 Data File : BM048810.D
 Acq On : 22 Nov 2024 10:58
 Operator : RC/JU
 Sample : SST001061
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Nov 22 14:29:51 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 22 14:18:42 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
36) 2-Methylnaphthalene	12.063	142	307733	9.823	ng/ul	98
37) 1-Methylnaphthalene	12.280	142	313575	9.827	ng/ul	99
39) 1,2,4,5-Tetrachloroben...	12.433	216	145779	9.042	ng/ul	99
40) Hexachlorocyclopentadiene	12.422	237	50497	6.197	ng/ul	93
41) 2,4,6-Trichlorophenol	12.680	196	61248	7.372	ng/ul	95
42) 2,4,5-Trichlorophenol	12.745	196	73868	8.133	ng/ul	96
43) 1,1'-Biphenyl	13.086	154	387153	9.468	ng/ul	99
44) 2-Chloronaphthalene	13.122	162	300846	9.511	ng/ul	100
45) 2-Nitroaniline	13.327	65	55201	7.605	ng/ul	90
47) Dimethylphthalate	13.727	163	345408	9.496	ng/ul	99
48) 2,6-Dinitrotoluene	13.833	165	44027	7.532	ng/ul	97
50) Acenaphthylene	13.974	152	481068	9.799	ng/ul	100
51) 3-Nitroaniline	14.163	138	50342	7.647	ng/ul#	96
52) Acenaphthene	14.321	153	340049	9.814	ng/ul	99
53) 2,4-Dinitrophenol	14.369	184	4156	3.016	ng/ul#	71
55) 4-Nitrophenol	14.474	109	30927	6.715	ng/ul	86
56) Dibenzofuran	14.657	168	445298	9.864	ng/ul	98
57) 2,4-Dinitrotoluene	14.621	165	61246	7.716	ng/ul	98
58) 2,3,4,6-Tetrachlorophenol	14.886	232	51921	7.474	ng/ul	96
59) Diethylphthalate	15.104	149	354065	9.513	ng/ul	99
61) Fluorene	15.310	166	367051	9.872	ng/ul	99
62) 4-Chlorophenyl-phenyle...	15.316	204	170526	9.614	ng/ul	99
63) 4-Nitroaniline	15.321	138	57973	9.460	ng/ul	97
66) 4,6-Dinitro-2-methylph...	15.380	198	13904	4.220	ng/ul#	81
67) N-Nitrosodiphenylamine	15.527	169	294047	9.456	ng/ul	99
68) 4-Bromophenyl-phenylether	16.210	248	95736	8.999	ng/ul	99
69) Hexachlorobenzene	16.310	284	115657	9.438	ng/ul	96
70) Atrazine	16.492	200	82186	8.077	ng/ul	99
71) Pentachlorophenol	16.663	266	30978	4.468	ng/ul	94
72) Phenanthrene	17.045	178	548831	9.754	ng/ul	99
74) Anthracene	17.139	178	551963	9.714	ng/ul	99
75) 1,2,3,4-Tetrachloroben...	13.045	216	142079	8.796	ng/uL	97
76) Pentachlorobenzene	14.574	250	141467	9.139	ng/uL	98
77) Carbazole	17.410	167	484509	9.528	ng/ul	99
78) Di-n-butylphthalate	18.009	149	485922	8.046	ng/ul	99
80) Fluoranthene	19.074	202	551248	9.043	ng/ul	99
82) Pyrene	19.433	202	618779	9.417	ng/ul	99
83) Butylbenzylphthalate	20.374	149	105951	5.609	ng/ul	90
84) 3,3'-Dichlorobenzidine	21.168	252	107369	6.947	ng/ul	97
85) Benzo(a)anthracene	21.227	228	548355	9.291	ng/ul	98
86) Bis(2-ethylhexyl)phtha...	21.209	149	210626	6.479	ng/ul	100
87) Chrysene	21.286	228	542738	9.723	ng/ul	98
89) Di-n-octyl phthalate	22.315	149	267489	4.761	ng/ul	100
90) Benzo(b)fluoranthene	23.209	252	484004	8.970	ng/ul	99
91) Benzo(k)fluoranthene	23.268	252	531455	9.312	ng/ul	98
93) Benzo(a)pyrene	23.968	252	474977	9.168	ng/ul	99
94) Indeno(1,2,3-cd)pyrene	27.291	276	557864	9.004	ng/ul	98
95) Dibenzo(a,h)anthracene	27.356	278	445473	8.906	ng/ul	99
96) Benzo(g,h,i)perylene	28.274	276	440715	9.148	ng/ul	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

[illegible]