

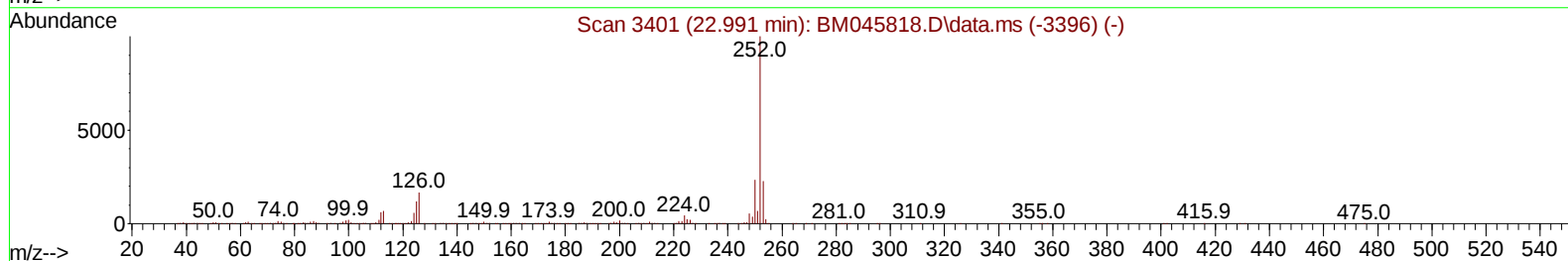
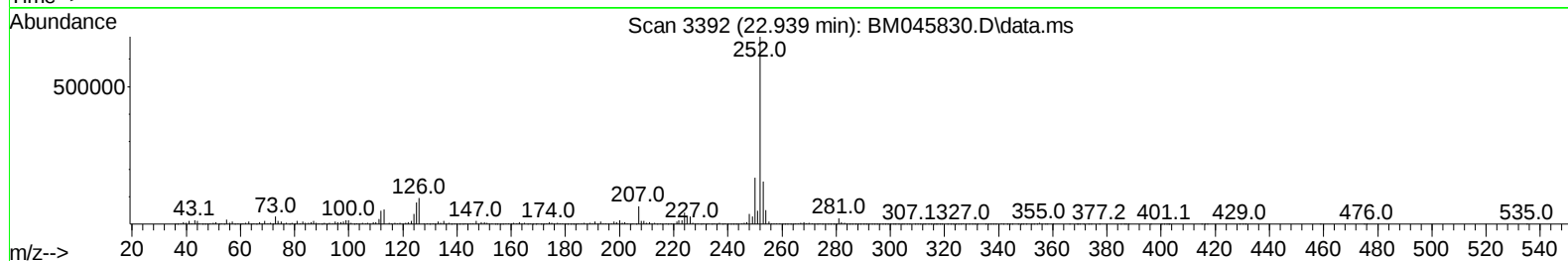
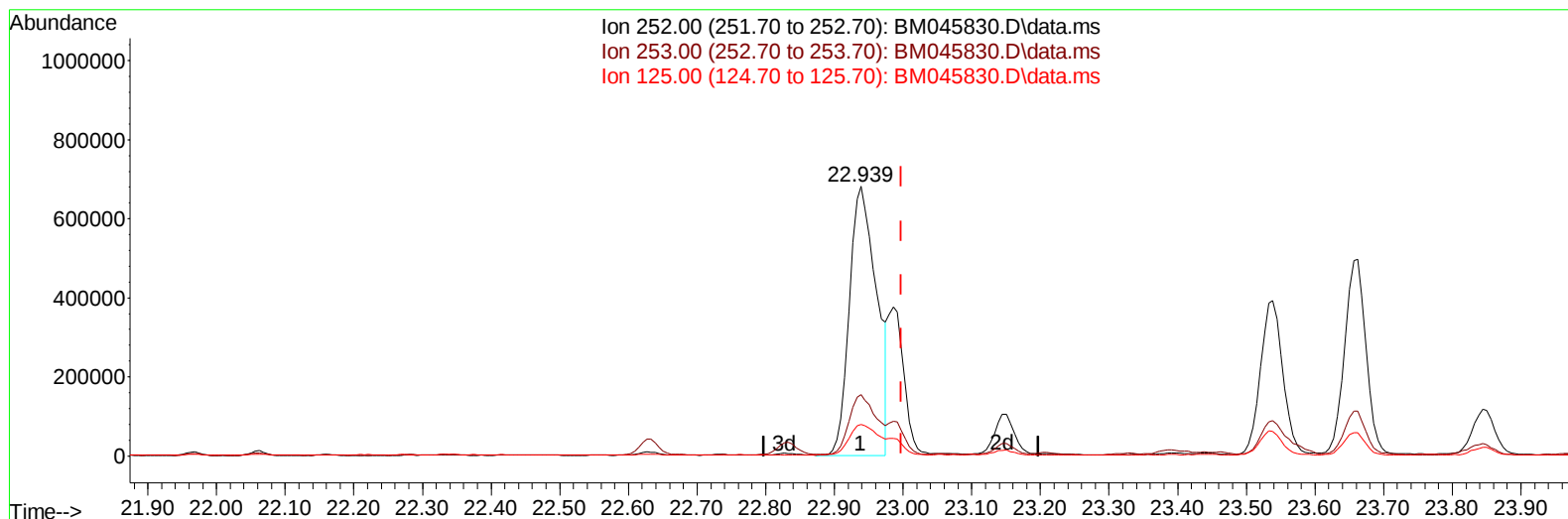
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060424\
 Data File : BM045830.D
 Acq On : 04 Jun 2024 23:50
 Operator : MA/JU
 Sample : P2589-19
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BH6D3

Manual IntegrationsAPPROVED

Quant Time: Jun 05 00:27:20 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM053124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 31 13:01:11 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/05/2024
 Supervised By :mohammad ahmed 06/07/2024



TIC: BM045830.D\data.ms

(91) Benzo(k)fluoranthene

22.939min (-0.059) 31.47 ng/u1

response 1876090

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	21.00	22.66
125.00	10.30	11.65
0.00	0.00	0.00

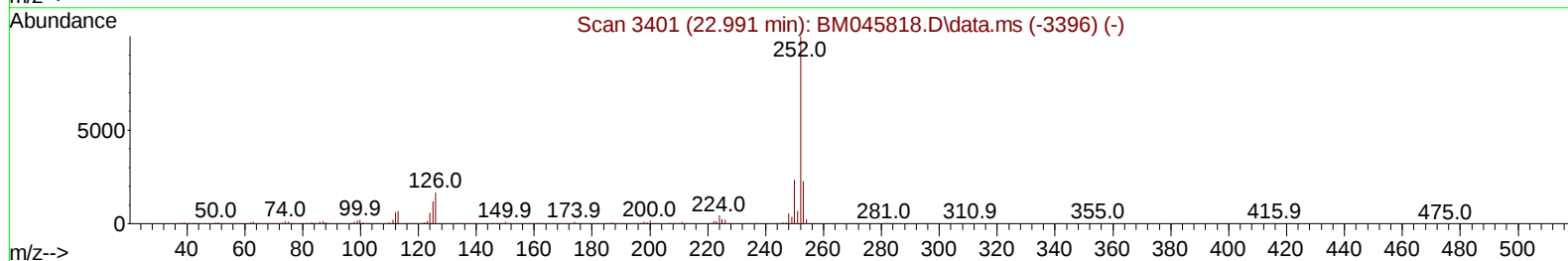
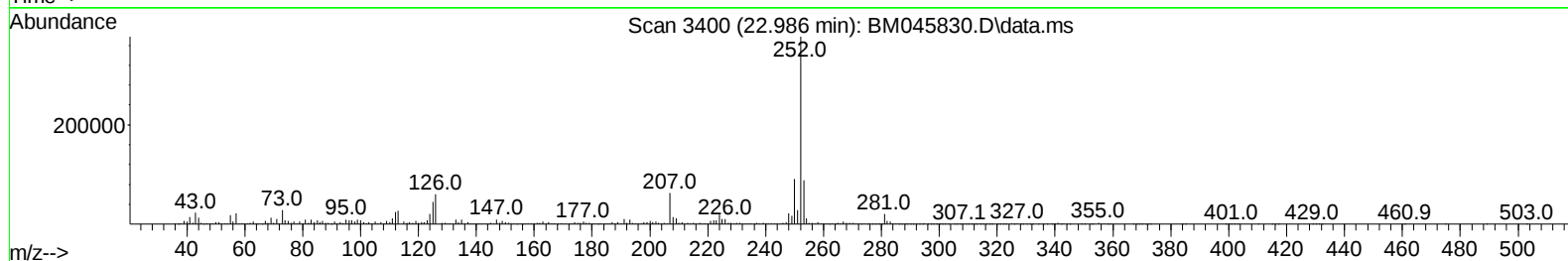
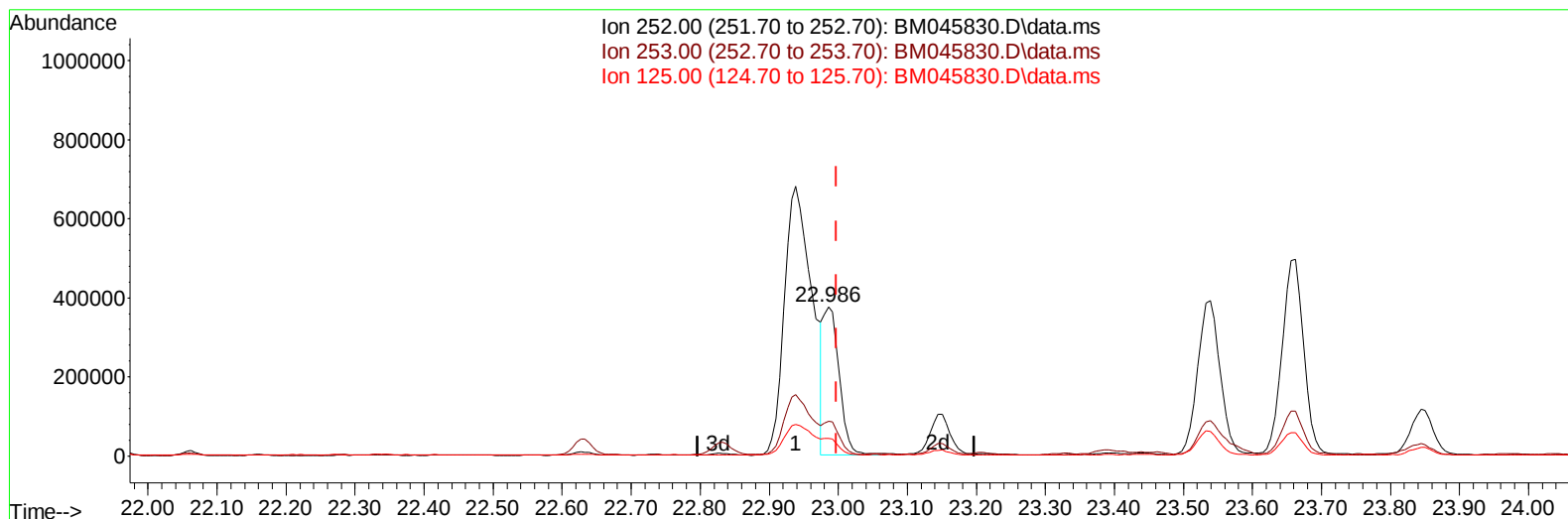
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TIC: BM045830.D\data.ms

(91) Benzo(k)fluoranthene

22.986min (-0.012) 10.04 ng/u1 m

response 598410

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	21.00	23.39
125.00	10.30	11.81
0.00	0.00	0.00

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 Sample : P2589-19
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

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Quant Time: Jun 05 00:28:25 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM053124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 31 13:01:11 2024
 Response via : Initial Calibration

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.463	152		305141	20.000	ng/ul	0.00
20) Naphthalene-d8	10.234	136		1296619	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.110	164		812543	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.851	188		1628509	20.000	ng/ul	0.00
79) Chrysene-d12	21.062	240		1043856	20.000	ng/ul	0.00
88) Perylene-d12	23.786	264		993030	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.052	96		25854	3.285	ng/uL	0.00
4) Pyridine-d5	3.469	84		228886	11.295	ng/ul	0.00
7) Phenol-d5	6.722	99		589369	22.736	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.828	67		430208	23.112	ng/ul	0.00
11) 2-Chlorophenol-d4	7.028	132		483556	24.973	ng/ul	0.00
15) 4-Methylphenol-d8	8.234	113		406589	20.761	ng/ul	0.00
21) Nitrobenzene-d5	8.628	128		245533	24.220	ng/ul	0.00
24) 2-Nitrophenol-d4	9.340	143		264294	25.759	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.892	165		493908	25.670	ng/ul	0.00
31) 4-Chloroaniline-d4	10.416	131		369403	13.872	ng/ul	0.00
46) Dimethylphthalate-d6	13.557	166		1555981	27.661	ng/ul	0.00
49) Acenaphthylene-d8	13.798	160		1686273	25.815	ng/ul	0.00
54) 4-Nitrophenol-d4	14.404	143		207844	17.120	ng/ul	0.00
60) Fluorene-d10	15.110	176		1316151	27.426	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.245	200		101814	12.243	ng/ul	0.00
73) Anthracene-d10	16.951	188		2004546	28.315	ng/ul	0.00
81) Pyrene-d10	19.251	212		2189157	33.723	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.603	264		1353487	29.163	ng/ul	0.00
Target Compounds							
						Qvalue	
50) Acenaphthylene	13.827	152		127605	1.688	ng/ul	99
52) Acenaphthene	14.175	153		88593	1.756	ng/ul	98
56) Dibenzofuran	14.510	168		107865	1.585	ng/ul	99
61) Fluorene	15.163	166		103305	1.885	ng/ul#	96
72) Phenanthrene	16.898	178		2446655	28.513	ng/ul	100
74) Anthracene	16.986	178		417401	4.870	ng/ul	98
77) Carbazole	17.274	167		292365	3.998	ng/ul	99
78) Di-n-butylphthalate	17.845	149		362105	3.765	ng/ul	99
80) Fluoranthene	18.915	202		4420233	56.167	ng/ul	98
82) Pyrene	19.280	202		3703321	45.801	ng/ul	98
83) Butylbenzylphthalate	20.209	149		292222	8.868	ng/ul	98
85) Benzo(a)anthracene	21.045	228		1339760	18.588	ng/ul	96
86) Bis(2-ethylhexyl)phtha...	20.998	149		1548807	31.634	ng/ul	100
87) Chrysene	21.103	228		1569940	23.546	ng/ul	98
90) Benzo(b)fluoranthene	22.939	252		1876090	30.742	ng/ul	98
91) Benzo(k)fluoranthene	22.986	252		598410m	10.039	ng/ul	
93) Benzo(a)pyrene	23.662	252		1071335	19.978	ng/ul	99
94) Indeno(1,2,3-cd)pyrene	26.785	276		809459	13.822	ng/ul	100
95) Dibenzo(a,h)anthracene	26.821	278		193239	4.172	ng/ul#	93
96) Benzo(g,h,i)perylene	27.715	276		704222	13.653	ng/ul	99

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

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