

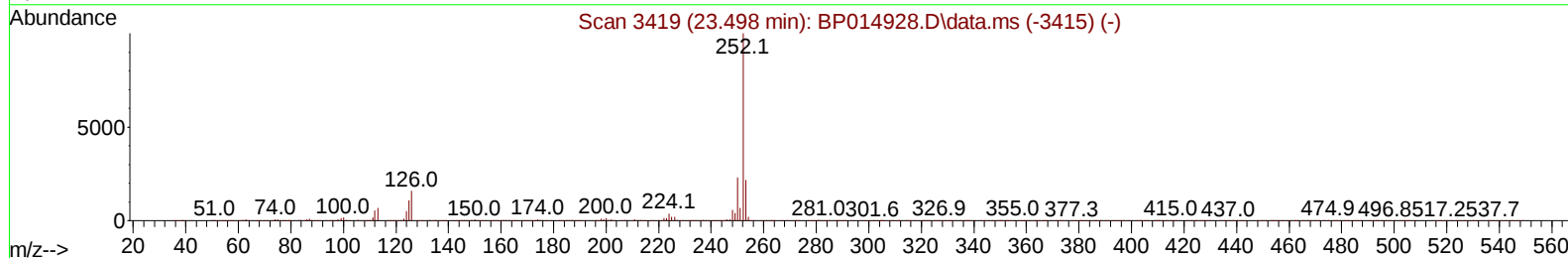
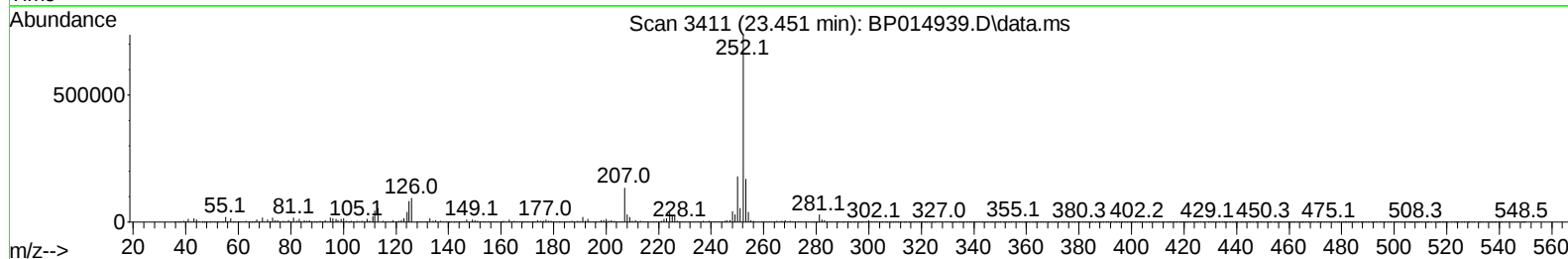
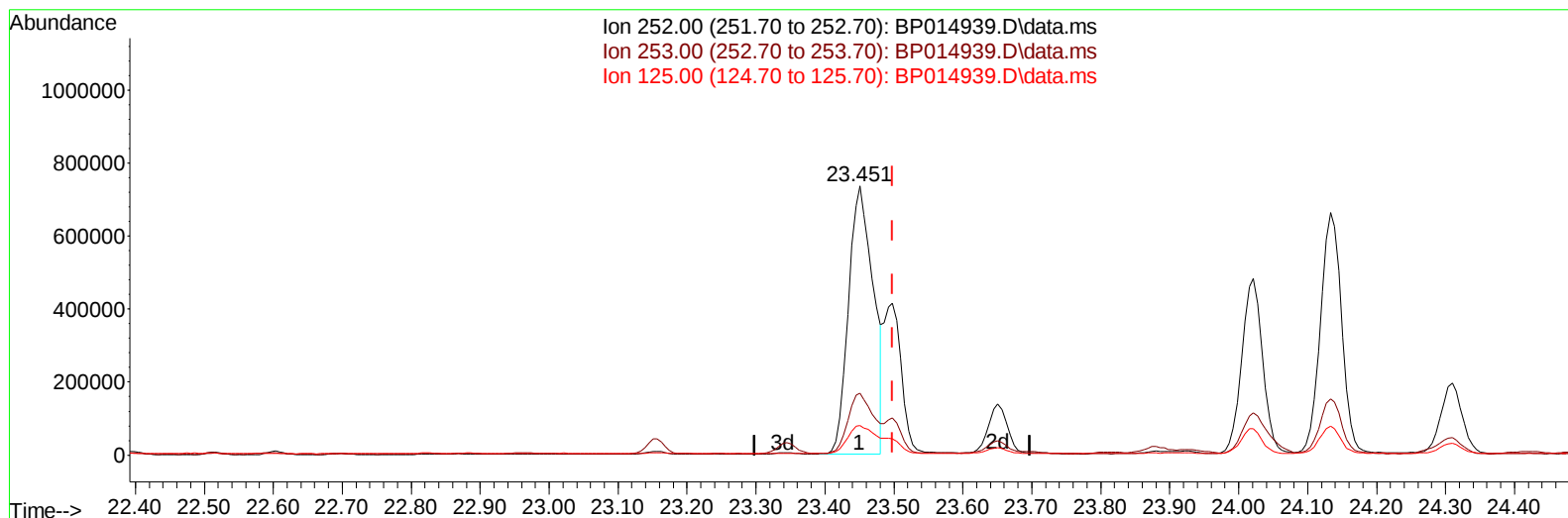
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050323\
 Data File : BP014939.D
 Acq On : 03 May 2023 20:17
 Operator : CG/JU
 Sample : 02419-22DL 2X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EW919DL

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 05/04/2023
 Supervised By :Jagrut Upadhyay 05/04/2023

Quant Time: May 04 09:10:58 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 03 23:33:17 2023
 Response via : Initial Calibration



TIC: BP014939.D\data.ms

(91) Benzo(k)fluoranthene

23.451min (-0.047) 24.47 ng/u1

response 1834161

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	22.20	22.90
125.00	11.10	11.03
0.00	0.00	0.00

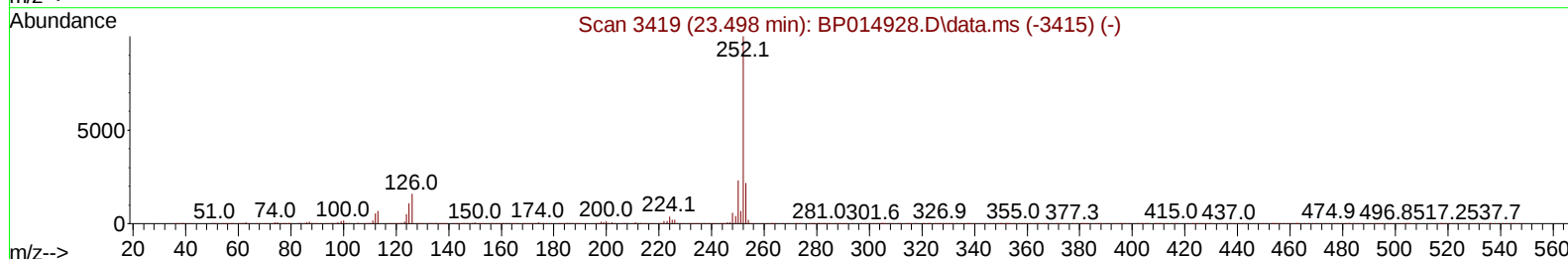
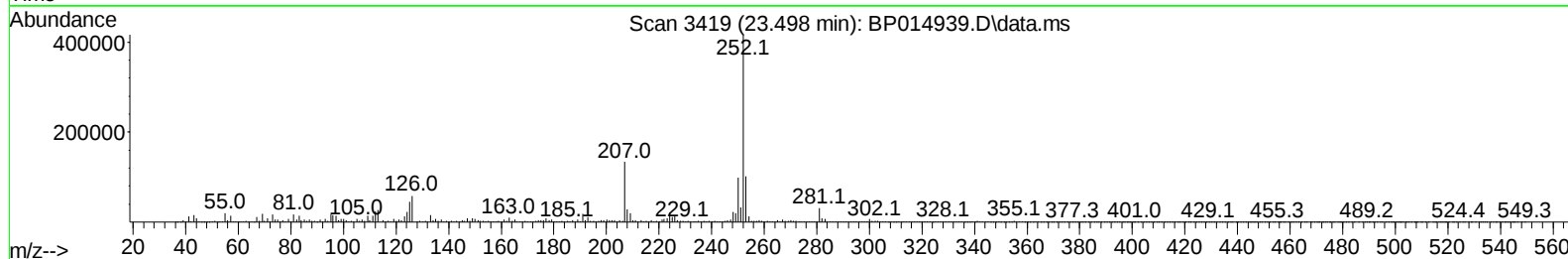
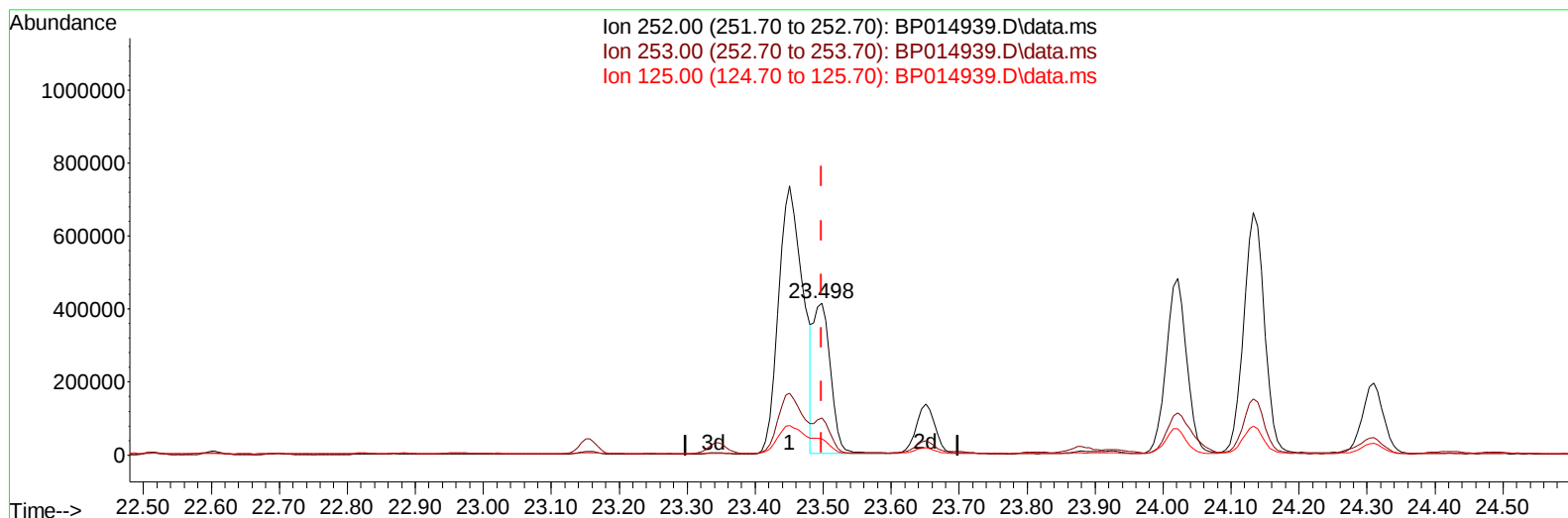
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(91) Benzo(k)fluoranthene

23.498min (-0.000) 9.75 ng/ul m

response 730971

Ion	Exp%	Act%
252.00	100.00	100.00
253.00	22.20	24.36
125.00	11.10	10.85
0.00	0.00	0.00

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

Internal Standards							
1) 1,4-Dichlorobenzene-d4	8.163	152		314224	20.000	ng/ul	0.00
20) Naphthalene-d8	10.998	136		1299834	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.810	164		686597	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.569	188		1113364	20.000	ng/ul	0.00
79) Chrysene-d12	21.651	240		887151	20.000	ng/ul	0.00
88) Perylene-d12	24.251	264		1139548	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.464	96		11255	1.332	ng/uL	0.00
4) Pyridine-d5	0.000	84		0d	0.000	ng/ul	
7) Phenol-d5	7.305	99		204476	7.506	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.481	67		133375	7.921	ng/ul	0.00
11) 2-Chlorophenol-d4	7.687	132		167555	8.365	ng/ul	0.00
15) 4-Methylphenol-d8	8.869	113		152453	7.262	ng/ul	0.00
21) Nitrobenzene-d5	9.340	128		74450	8.053	ng/ul	0.00
24) 2-Nitrophenol-d4	10.069	143		75987	8.005	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.610	165		145577	7.810	ng/ul	0.00
31) 4-Chloroaniline-d4	11.134	131		82852	3.073	ng/ul	0.00
46) Dimethylphthalate-d6	14.210	166		445949	9.177	ng/ul	0.00
49) Acenaphthylene-d8	14.504	160		510479	8.759	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143		0d	0.000	ng/ul	
60) Fluorene-d10	15.798	176		356013	8.516	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.910	200		6955	1.110	ng/ul	0.00
73) Anthracene-d10	17.663	188		457436	9.144	ng/ul	0.00
81) Pyrene-d10	19.880	212		472233	7.892	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.080	264		536917	9.670	ng/ul	0.00
Target Compounds							
						Qvalue	
36) 2-Methylnaphthalene	12.645	142		47614	1.078	ng/ul	100
52) Acenaphthene	14.875	153		95935	2.098	ng/ul	98
56) Dibenzofuran	15.204	168		80733	1.309	ng/ul	99
61) Fluorene	15.857	166		96787	1.989	ng/ul	97
72) Phenanthrene	17.610	178		1649649	26.820	ng/ul	100
74) Anthracene	17.698	178		395107	6.452	ng/ul	99
77) Carbazole	17.963	167		146067	2.836	ng/ul	99
80) Fluoranthene	19.557	202		2785271	37.936	ng/ul	99
82) Pyrene	19.910	202		2481599	31.902	ng/ul	99
85) Benzo(a)anthracene	21.633	228		1274405	21.522	ng/ul	99
87) Chrysene	21.686	228		1254539	21.041	ng/ul	98
90) Benzo(b)fluoranthene	23.451	252		1834161	24.905	ng/ul	98
91) Benzo(k)fluoranthene	23.498	252		730971m	9.752	ng/ul	
93) Benzo(a)pyrene	24.133	252		1395859	21.612	ng/ul	98
94) Indeno(1,2,3-cd)pyrene	27.003	276		1052465	13.359	ng/ul	96
95) Dibenzo(a,h)anthracene	27.015	278		264593	4.098	ng/ul#	81
96) Benzo(g,h,i)perylene	27.862	276		979938	14.614	ng/ul	98

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

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