

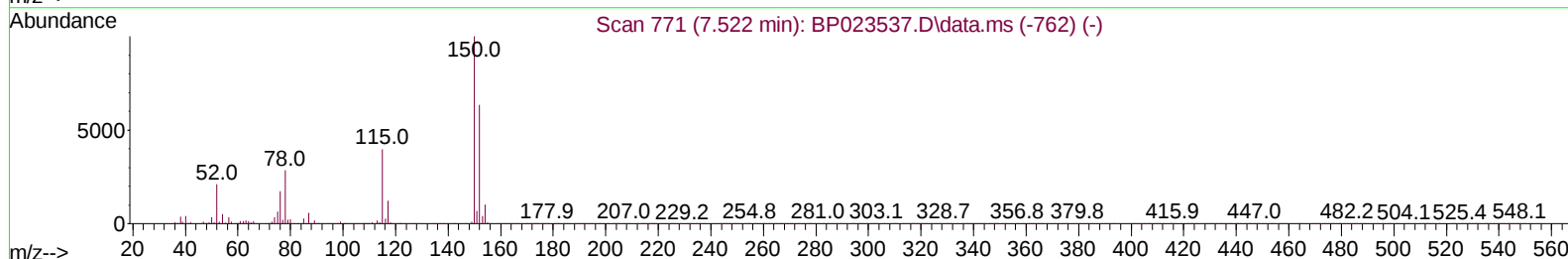
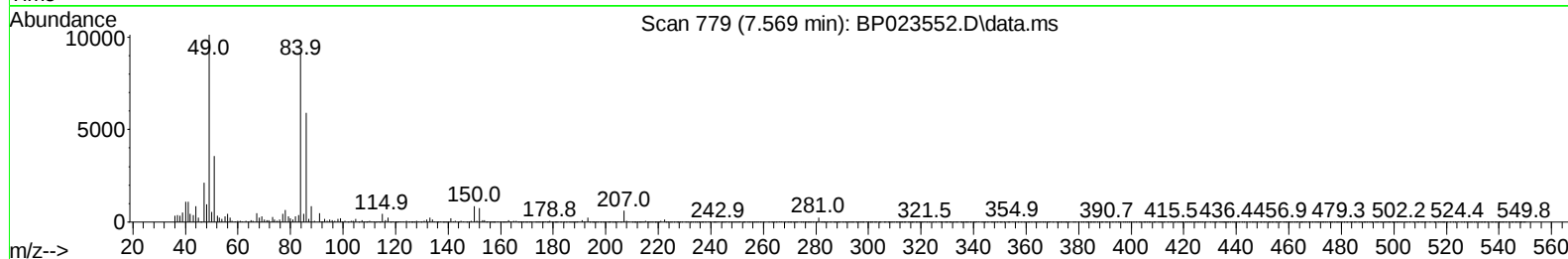
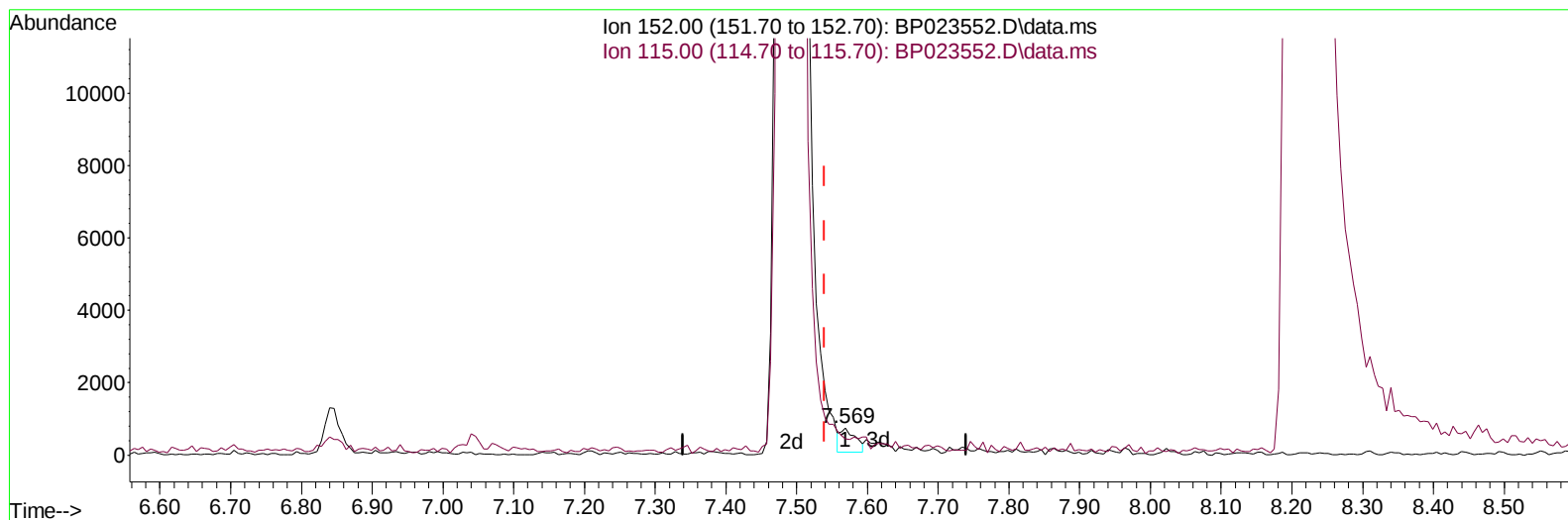
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121824\
 Data File : BP023552.D
 Acq On : 20 Dec 2024 10:32
 Operator : RC/JU
 Sample : PB165707BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK707

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 12/21/2024
 Supervised By :mohammad ahmed 12/21/2024

Quant Time: Dec 20 23:05:04 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 11 00:30:30 2024
 Response via : Initial Calibration



TIC: BP023552.D\data.ms

(1) 1,4-Dichlorobenzene-d4 (I)

7.569min (+ 0.029) 20.00 ng/u1

response 971

Ion	Exp%	Act%
152.00	100.00	100.00
115.00	65.20	59.09
0.00	0.00	0.00
0.00	0.00	0.00

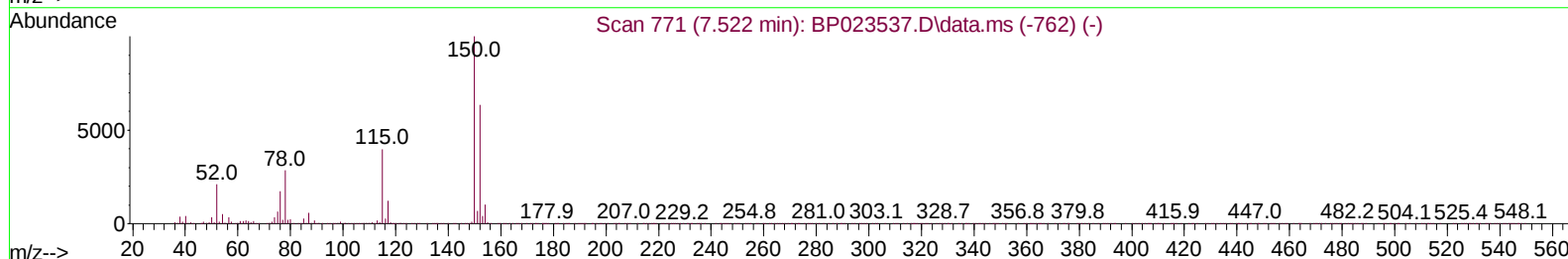
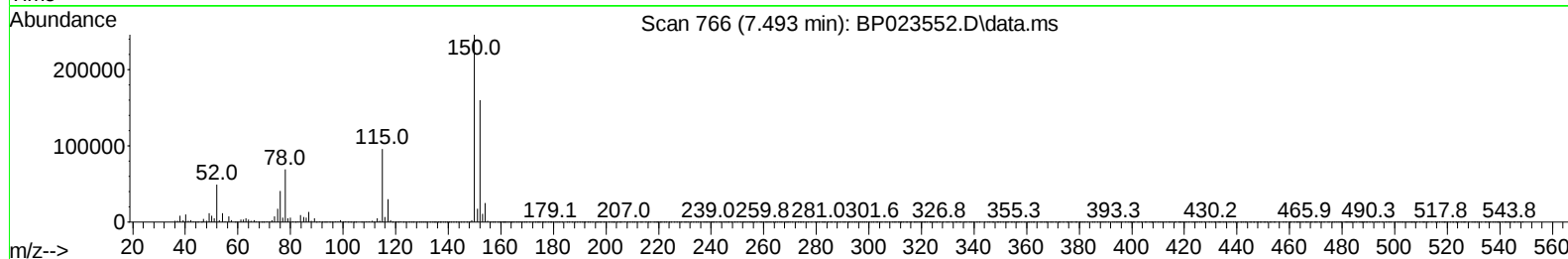
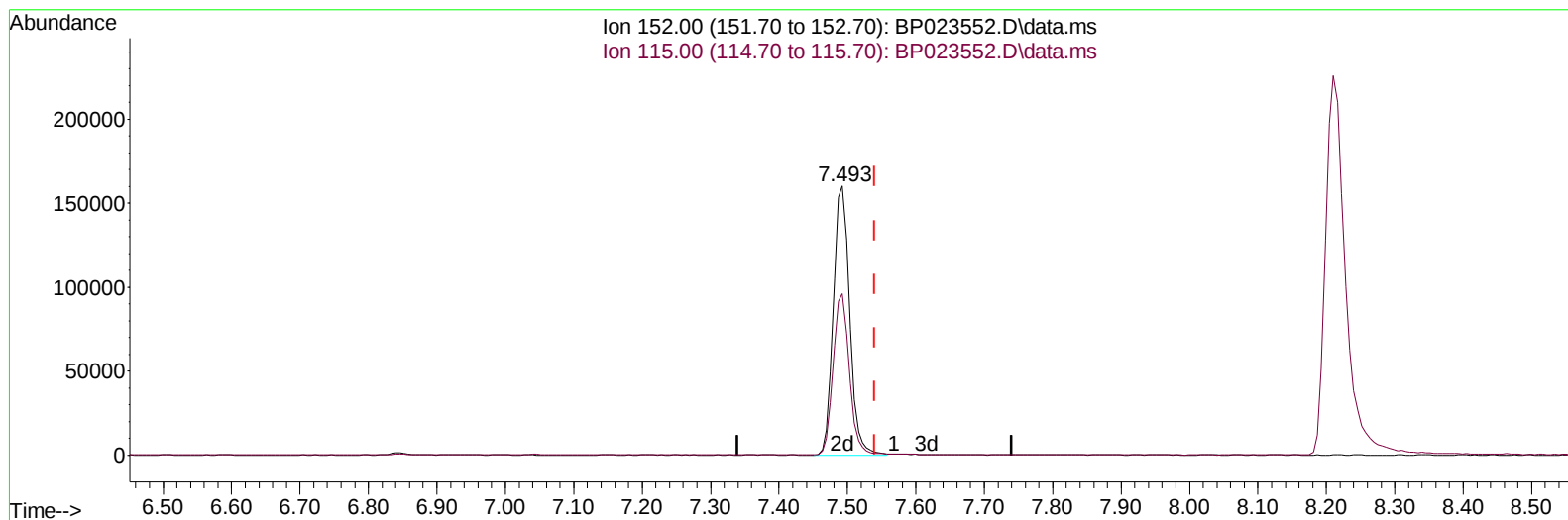
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TIC: BP023552.D\data.ms

(1) 1,4-Dichlorobenzene-d4 (I)

7.493min (-0.047) 20.00 ng/u1 m

response 266833

Ion	Exp%	Act%
152.00	100.00	100.00
115.00	65.20	59.94
0.00	0.00	0.00
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 12/21/2024
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Quant Time: Dec 20 23:28:48 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112624.MA.M
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.493	152		266833m	20.000	ng/ul	-0.05
20) Naphthalene-d8	10.240	136		961255	20.000	ng/ul	#-0.04
38) Acenaphthene-d10	14.122	164		700853	20.000	ng/ul	-0.02
64) Phenanthrene-d10	16.939	188		1591571	20.000	ng/ul	#-0.02
79) Chrysene-d12	21.380	240		1808894	20.000	ng/ul	# 0.00
88) Perylene-d12	24.545	264		1983526	20.000	ng/ul	-0.03
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.075	96		40177	5.543	ng/uL	-0.03
4) Pyridine-d5	3.481	84		529593	28.416	ng/ul	-0.04
7) Phenol-d5	6.705	99		661700	31.605	ng/ul	-0.04
9) Bis-(2-Chloroethyl)eth...	6.846	67		409092	30.971	ng/ul	-0.05
11) 2-Chlorophenol-d4	7.040	132		479804	31.051	ng/ul	-0.05
15) 4-Methylphenol-d8	8.210	113		503637	29.805	ng/ul	-0.05
21) Nitrobenzene-d5	8.634	128		227938	31.755	ng/ul	-0.05
24) 2-Nitrophenol-d4	9.351	143		276397	33.410	ng/ul	-0.04
28) 2,4-Dichlorophenol-d3	9.899	165		482174	27.616	ng/ul	-0.04
31) 4-Chloroaniline-d4	10.393	131		592397	30.286	ng/ul	-0.05
46) Dimethylphthalate-d6	13.539	166		1695438	31.162	ng/ul	-0.02
49) Acenaphthylene-d8	13.810	160		1790845	31.206	ng/ul	-0.02
54) 4-Nitrophenol-d4	14.434	143		240851	36.124	ng/ul	0.02
60) Fluorene-d10	15.145	176		1456978	30.445	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.310	200		302920	31.492	ng/ul	0.00
73) Anthracene-d10	17.039	188		2476729	34.097	ng/ul	-0.01
81) Pyrene-d10	19.439	212		3038803	32.154	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.327	264		3270034	30.796	ng/ul	-0.02

Target Compounds Qvalue

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

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