

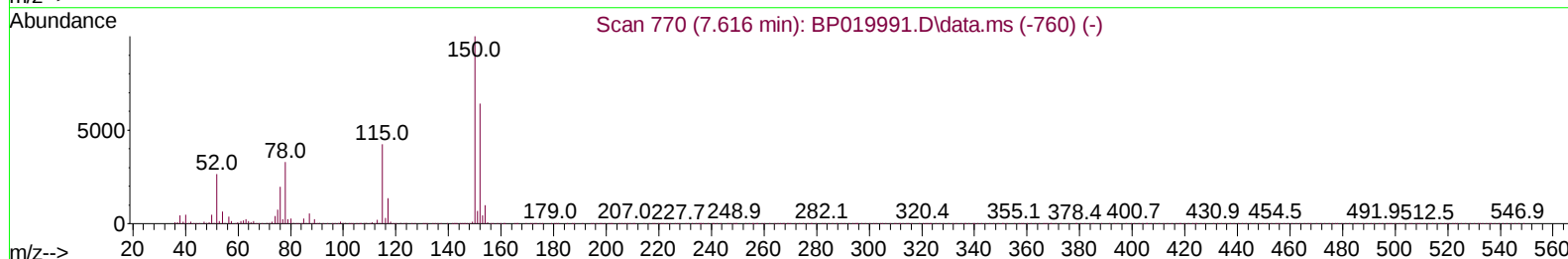
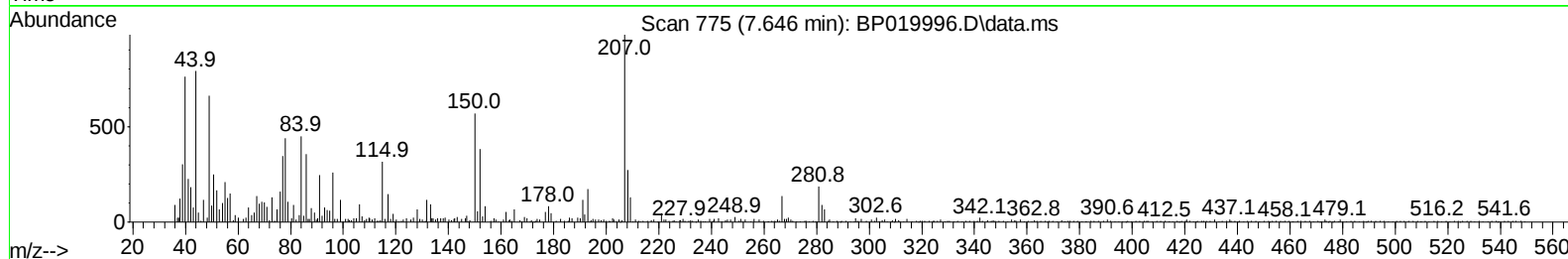
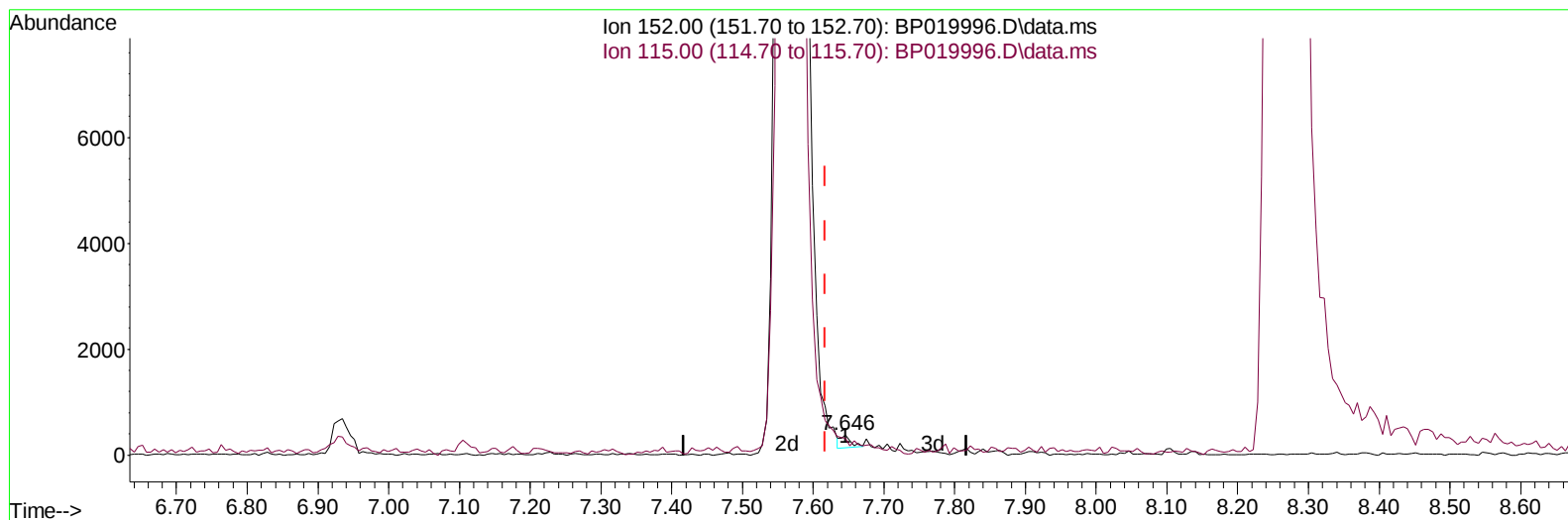
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032624\
 Data File : BP019996.D
 Acq On : 26 Mar 2024 14:02
 Operator : MA/JU
 Sample : PB159472BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK472

Manual IntegrationsAPPROVED

Quant Time: Mar 26 14:42:01 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Mar 25 10:46:59 2024
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 03/29/2024
 Supervised By :mohammad ahmed 03/29/2024



TIC: BP019996.D\data.ms

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

7.646min (+ 0.029) 20.00 ng/u1

response 243

Ion	Exp%	Act%
152.00	100.00	100.00
115.00	65.90	87.24#
0.00	0.00	0.00
0.00	0.00	0.00

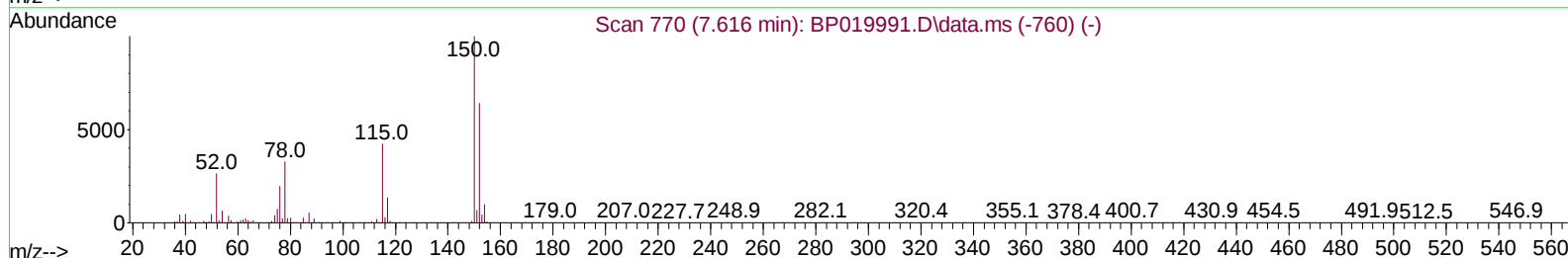
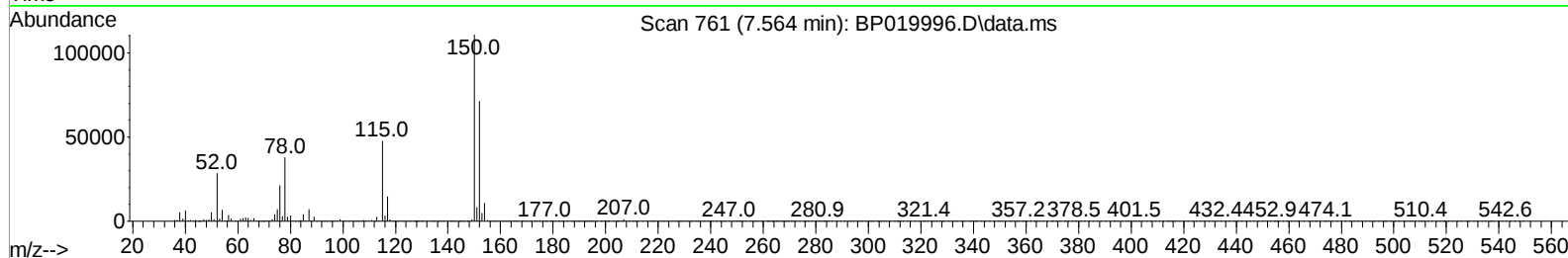
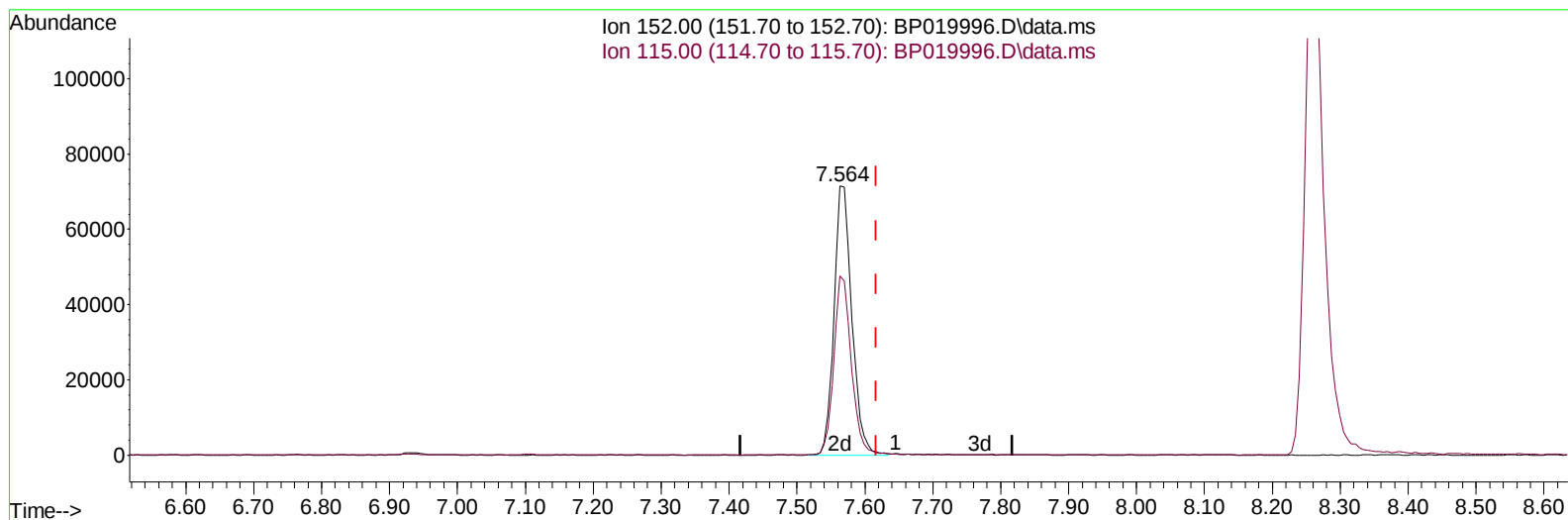
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TIC: BP019996.D\data.ms

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

7.564min (-0.053) 20.00 ng/ul m

response 129906

Ion	Exp%	Act%
152.00	100.00	100.00
115.00	65.90	66.56
0.00	0.00	0.00
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 03/29/2024
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Quant Time: Mar 26 14:43:19 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.564	152	129906m	20.000	ng/ul	-0.05
20) Naphthalene-d8	10.334	136	550195	20.000	ng/ul	-0.04
38) Acenaphthene-d10	14.234	164	387750	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.069	188	811237	20.000	ng/ul	0.00
79) Chrysene-d12	21.545	240	613858	20.000	ng/ul	0.00
88) Perylene-d12	24.833	264	610241	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	21535	7.263	ng/uL	-0.03
4) Pyridine-d5	3.611	84	308221	34.428	ng/ul	-0.03
7) Phenol-d5	6.752	99	377187	34.427	ng/ul	-0.05
9) Bis-(2-Chloroethyl)eth...	6.928	67	265407	37.182	ng/ul	-0.04
11) 2-Chlorophenol-d4	7.111	132	273030	34.700	ng/ul	-0.04
15) 4-Methylphenol-d8	8.258	113	304197	35.887	ng/ul	-0.05
21) Nitrobenzene-d5	8.722	128	148517	35.218	ng/ul	-0.04
24) 2-Nitrophenol-d4	9.434	143	168010	35.608	ng/ul	-0.04
28) 2,4-Dichlorophenol-d3	9.963	165	258620	30.012	ng/ul	-0.04
31) 4-Chloroaniline-d4	10.481	131	410566	36.548	ng/ul	-0.04
46) Dimethylphthalate-d6	13.651	166	1009468	37.132	ng/ul	0.00
49) Acenaphthylene-d8	13.922	160	1085518	34.181	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.475	143	129688	30.344	ng/ul	0.00
60) Fluorene-d10	15.263	176	842038	36.360	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.410	200	154217	31.758	ng/ul	-0.01
73) Anthracene-d10	17.175	188	1269799	35.360	ng/ul	0.00
81) Pyrene-d10	19.575	212	1433838	39.811	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.598	264	1113749	37.990	ng/ul	-0.01

Target Compounds Qvalue

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

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