

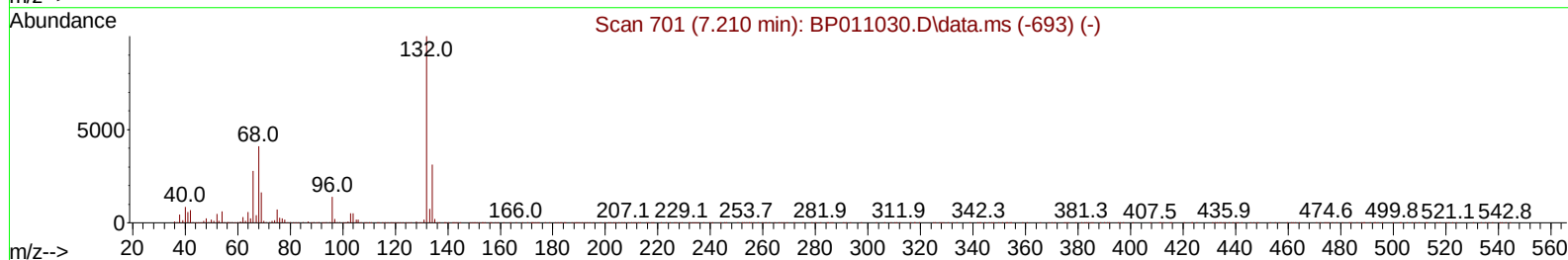
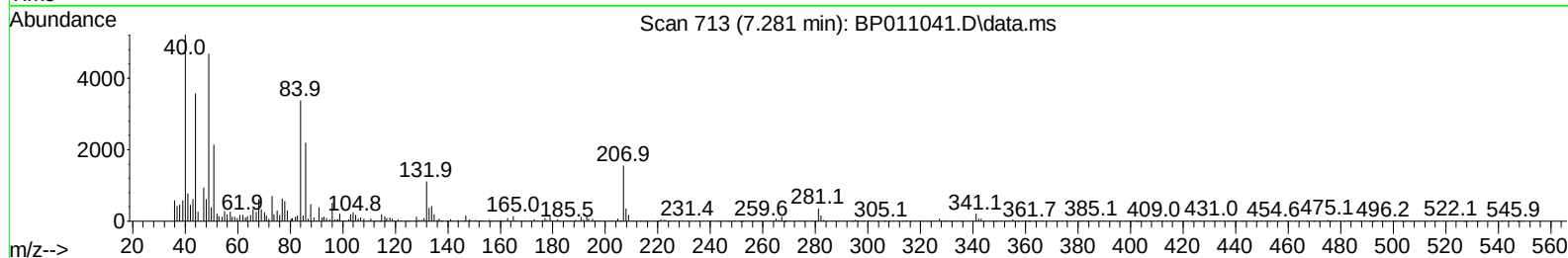
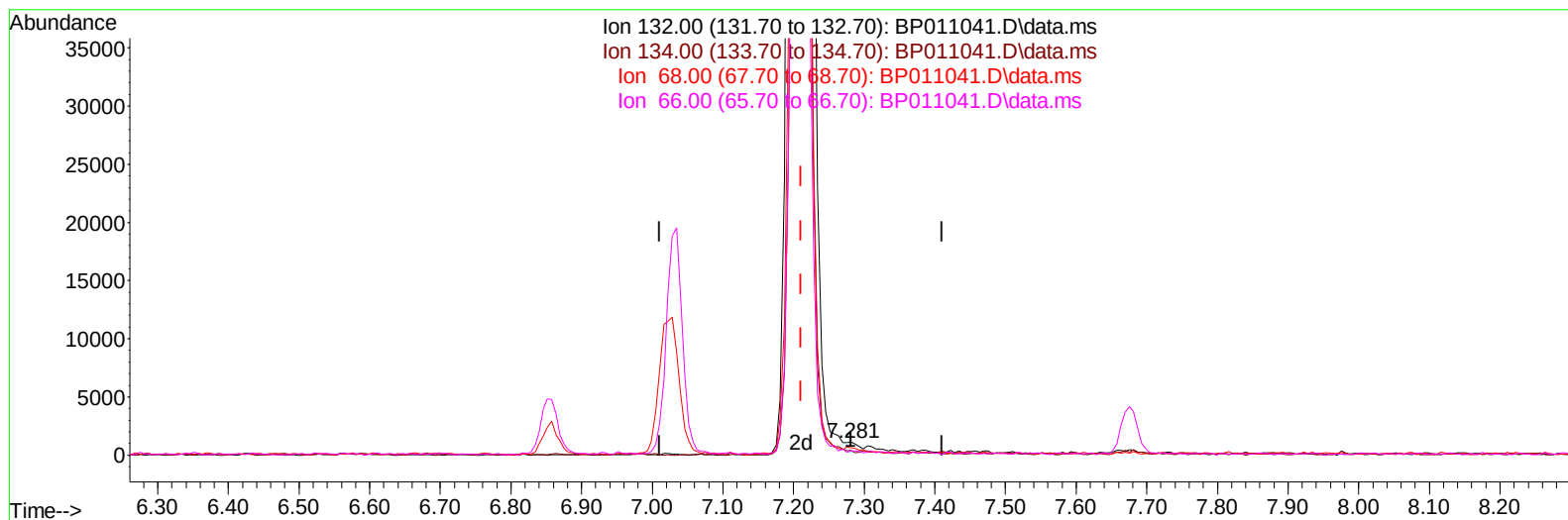
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071922\
 Data File : BP011041.D
 Acq On : 19 Jul 2022 20:09
 Operator : CG/JU
 Sample : N3753-09
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GBHQ0

Manual IntegrationsAPPROVED

Quant Time: Jul 19 23:41:45 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 19 03:25:54 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 07/20/2022
 Supervised By :mohammad ahmed 07/21/2022



TIC: BP011041.D\data.ms

(11) 2-Chlorophenol-d4 (S)

7.281min (+ 0.071) 0.04 ng/ul

response 771

Ion	Exp%	Act%
132.00	100.00	100.00
134.00	32.60	37.53
68.00	48.50	56.26
66.00	37.40	30.78

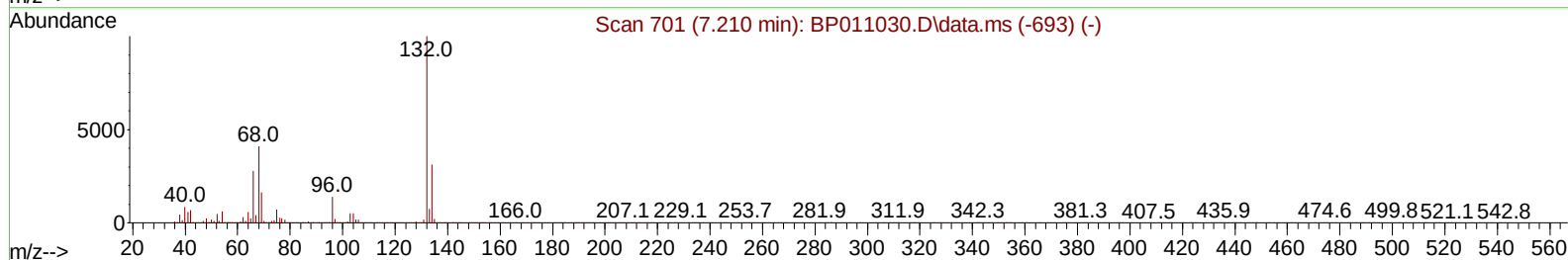
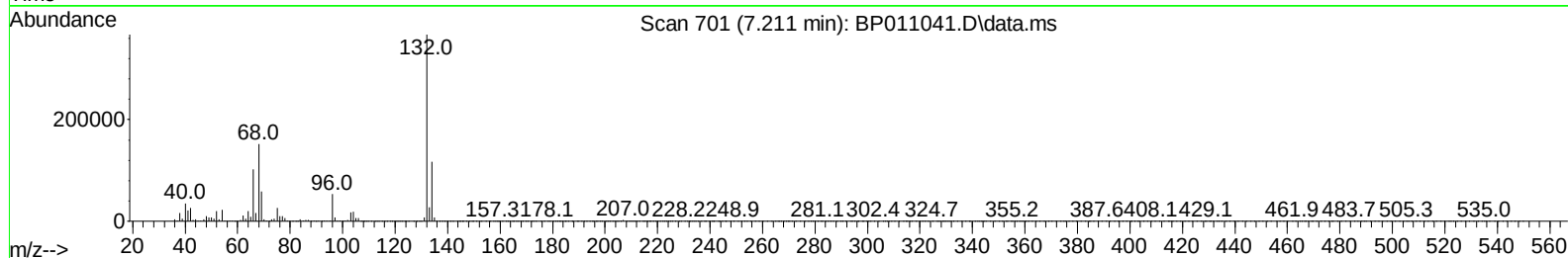
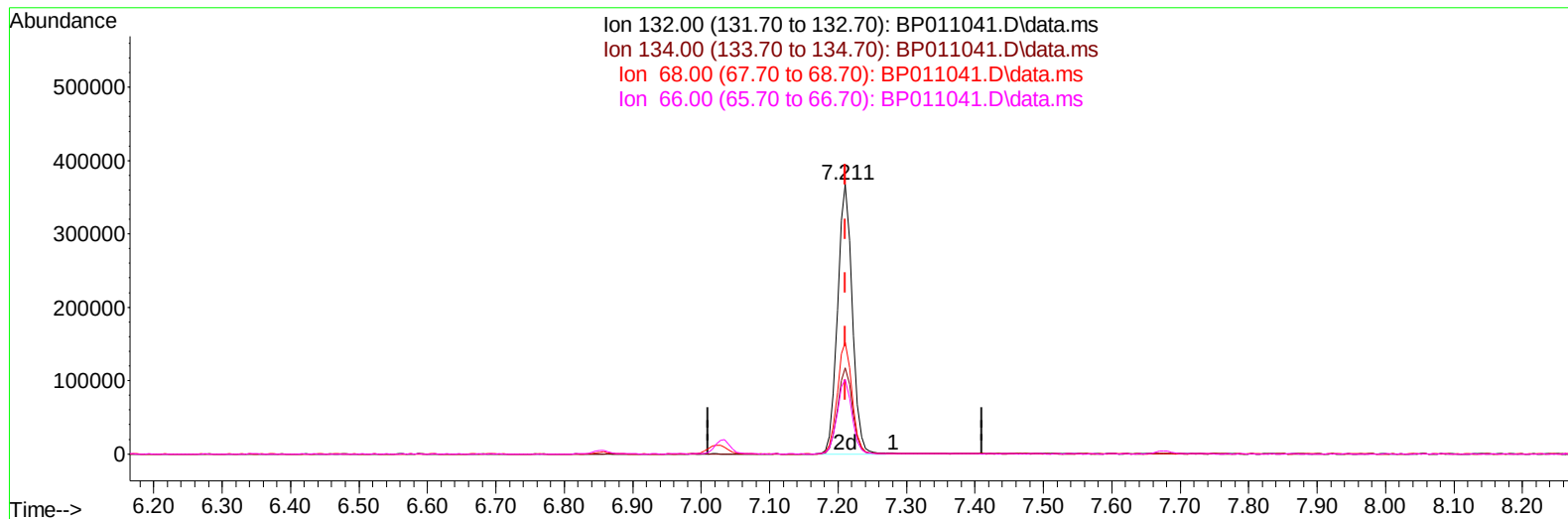
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(11) 2-Chlorophenol-d4 (S)

7.211min (+ 0.000) 24.98 ng/ul m

response 548494

Ion	Exp%	Act%
132.00	100.00	100.00
134.00	32.60	31.87
68.00	48.50	41.42
66.00	37.40	27.73#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.675	152	331099	20.000	ng/ul	0.00
20) Naphthalene-d8	10.469	136	1333783	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.340	164	740688	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.104	188	1479825	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.227	240	1448645	20.000	ng/ul	# 0.00
88) Perylene-d12	23.580	264	1477051	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	42707	4.650	ng/uL	0.00
4) Pyridine-d5	3.599	84	174934	7.369	ng/ul	0.00
7) Phenol-d5	6.852	99	160324	5.899	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.016	67	601825	34.436	ng/ul	0.00
11) 2-Chlorophenol-d4	7.211	132	548494m	24.975	ng/ul	0.00
15) 4-Methylphenol-d8	8.393	113	311756	14.328	ng/ul	0.00
21) Nitrobenzene-d5	8.840	128	344412	35.434	ng/ul	0.00
24) 2-Nitrophenol-d4	9.557	143	301386	31.336	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.093	165	501095	28.489	ng/ul	0.00
31) 4-Chloroaniline-d4	10.616	131	752149	25.939	ng/ul	0.00
46) Dimethylphthalate-d6	13.757	166	1913223	37.793	ng/ul	0.00
49) Acenaphthylene-d8	14.028	160	2255186	37.757	ng/ul	0.00
54) 4-Nitrophenol-d4	14.557	143	27593	2.693	ng/ul	0.00
60) Fluorene-d10	15.340	176	1633511	37.962	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.463	200	191391	26.137	ng/ul	0.00
73) Anthracene-d10	17.204	188	2549895	40.532	ng/ul	0.00
81) Pyrene-d10	19.463	212	2952640	38.808	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.427	264	2844068	40.090	ng/ul	0.00

Target Compounds Qvalue

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

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