

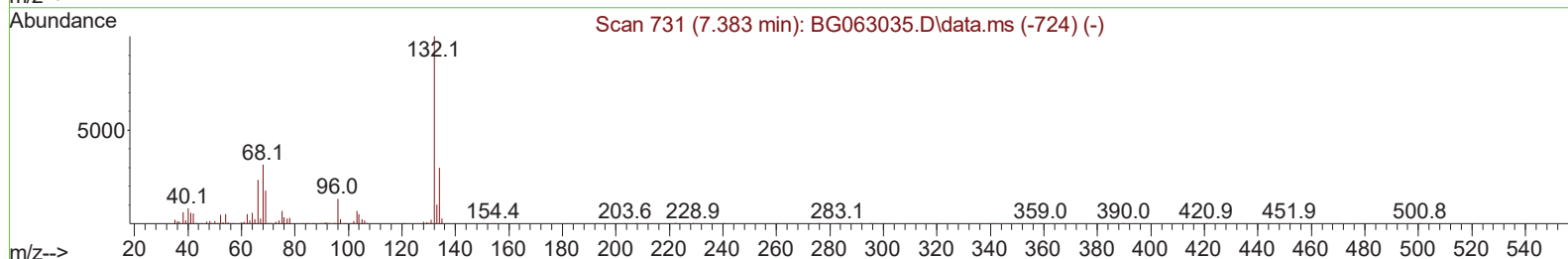
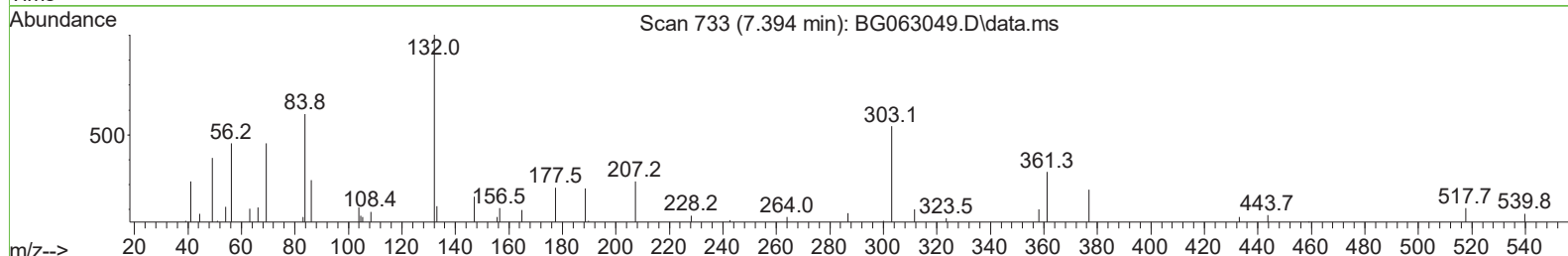
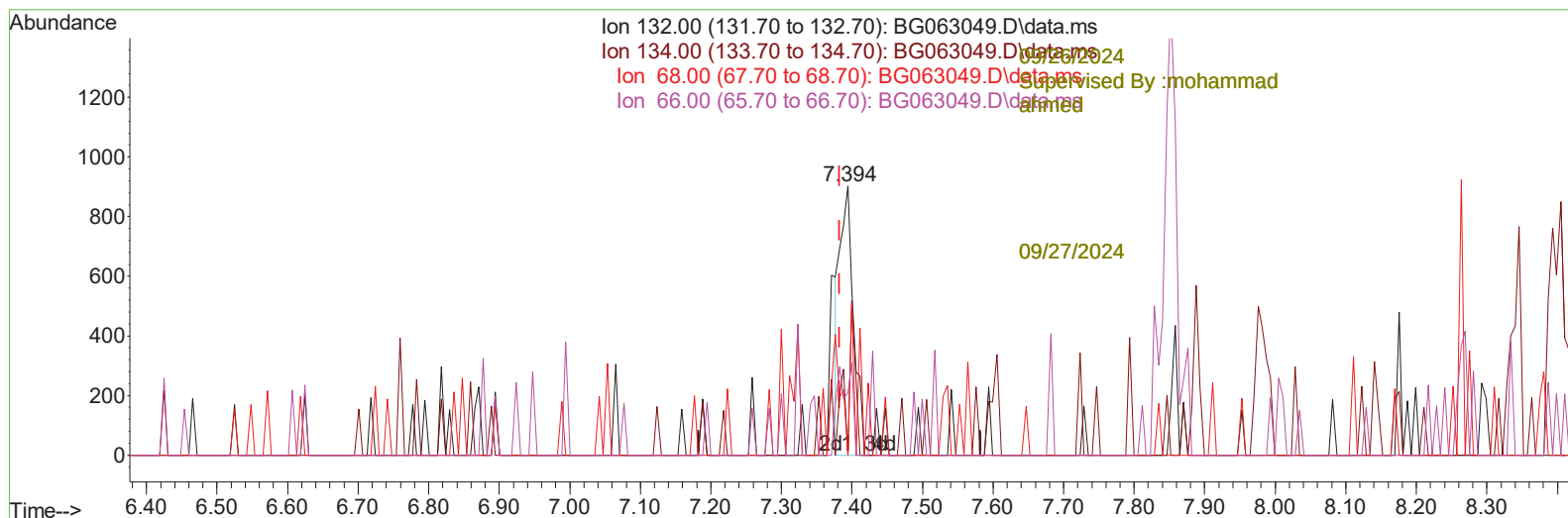
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092524\
Data File : BG063049.D
Acq On : 26 Sep 2024 1:17
Operator : RC/JU
Sample : P3879-17DL2 100X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DDC46DL2

Manual IntegrationsAPPROVED

Quant Time: Sep 26 01:58:15 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Sep 25 13:16:45 2024
Response via : Initial Calibration

Reviewed By :Yogesh
Patel



TIC: BG063049.D\data.ms

(11) 2-Chlorophenol-d4 (S)

7.394min (+ 0.011) 0.25 ng/uL

response 1221

Ion	Exp%	Act%
132.00	100.00	100.00
134.00	32.00	0.00#
68.00	25.00	0.00#
66.00	18.80	23.17#

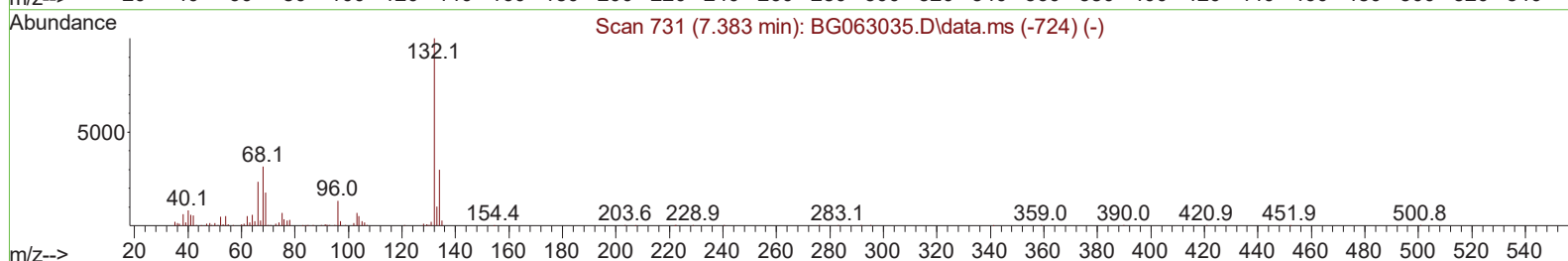
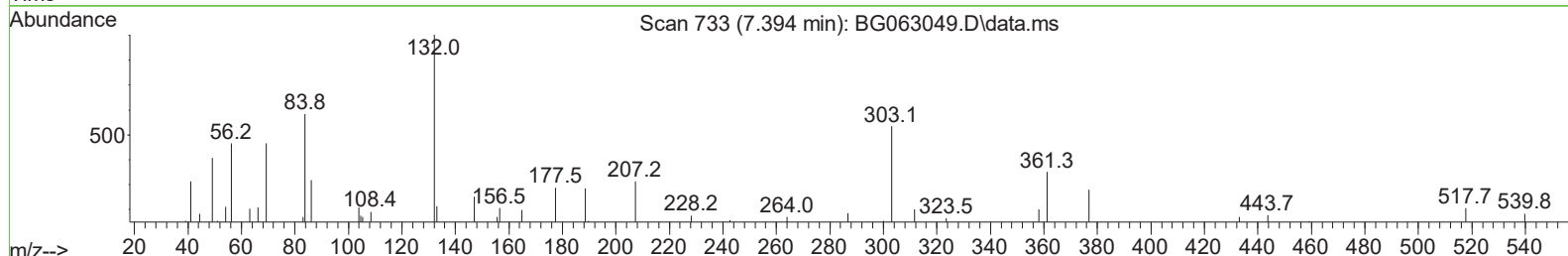
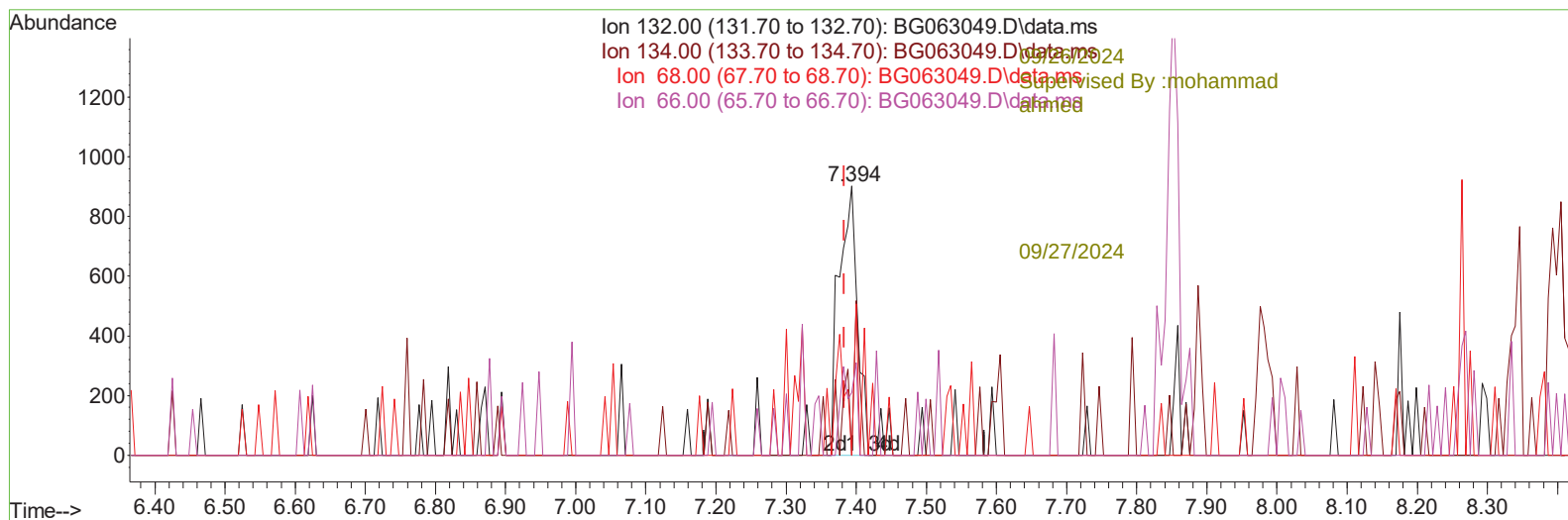
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TIC: BG063049.D\data.ms

(11) 2-Chlorophenol-d4 (S)

7.394min (+ 0.011) 0.33 ng/ul m

response 1644

Ion	Exp%	Act%
132.00	100.00	100.00
134.00	32.00	0.00#
68.00	25.00	0.00#
66.00	18.80	23.17#

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Manual IntegrationsAPPROVED

Quant Time: Sep 26 02:00:47 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG092524.MA.M
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 QLast Update : Wed Sep 25 13:16:45 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	82064	20.000	ng/ul	0.00
20) Naphthalene-d8	10.637	136	303343	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.486	164	263995	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.235	188	752986	20.000	ng/ul	0.00
79) Chrysene-d12	21.483	240	1081020	20.000	ng/ul	# 0.00
88) Perylene-d12	24.527	264	1261448	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.000	96	0	0.000	ng/uL	
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	0.000	99	0d	0.000	ng/ul	
9) Bis-(2-Chloroethyl)eth...	0.000	67	0d	0.000	ng/ul	
11) 2-Chlorophenol-d4	7.394	132	1644m	0.330	ng/ul	0.01
15) 4-Methylphenol-d8	0.000	113	0d	0.000	ng/ul	
21) Nitrobenzene-d5	0.000	128	0d	0.000	ng/ul	
24) 2-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
28) 2,4-Dichlorophenol-d3	0.000	165	0d	0.000	ng/ul	
31) 4-Chloroaniline-d4	0.000	131	0d	0.000	ng/ul	
46) Dimethylphthalate-d6	13.892	166	8739	0.416	ng/ul	0.00
49) Acenaphthylene-d8	14.168	160	7594	0.341	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.479	176	7718	0.401	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.626	200	5286	1.005	ng/ul	0.02
73) Anthracene-d10	17.335	188	17552	0.486	ng/ul	0.00
81) Pyrene-d10	19.627	212	20644	0.339	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.310	264	27827	0.412	ng/ul	0.00
Target Compounds						
58) 2,3,4,6-Tetrachlorophenol	15.114	232	98692	14.246	ng/ul#	95
71) Pentachlorophenol	16.889	266	265342	41.415	ng/ul	90

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

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