

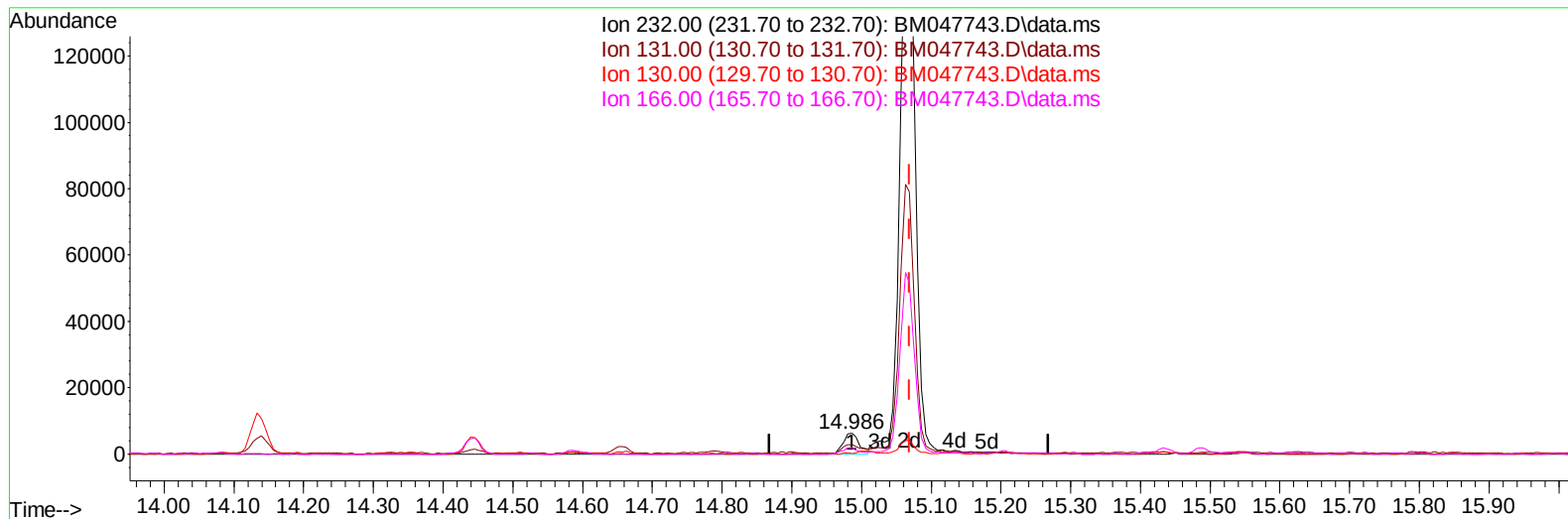
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047743.D
Acq On : 01 Oct 2024 15:03
Operator : RC/JU
Sample : P4018-06
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDCB2

Manual IntegrationsAPPROVED

Quant Time: Oct 01 22:49:16 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Sep 28 02:30:36 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/02/2024
Supervised By :mohammad ahmed 10/04/2024



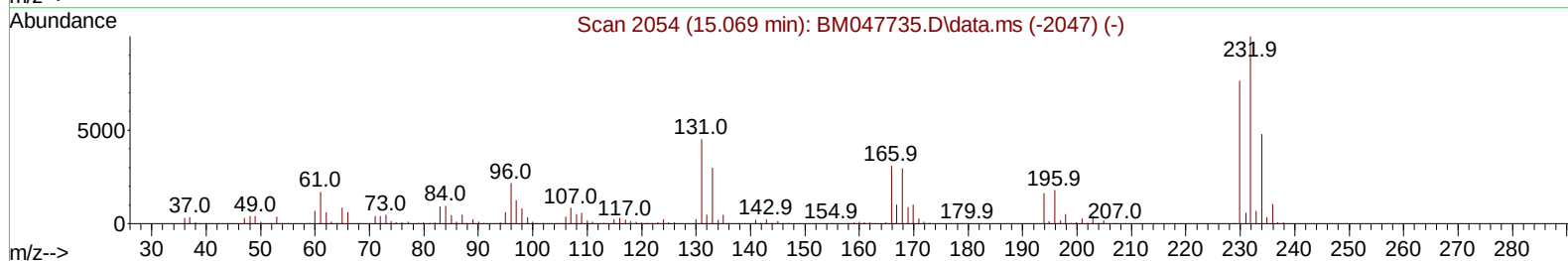
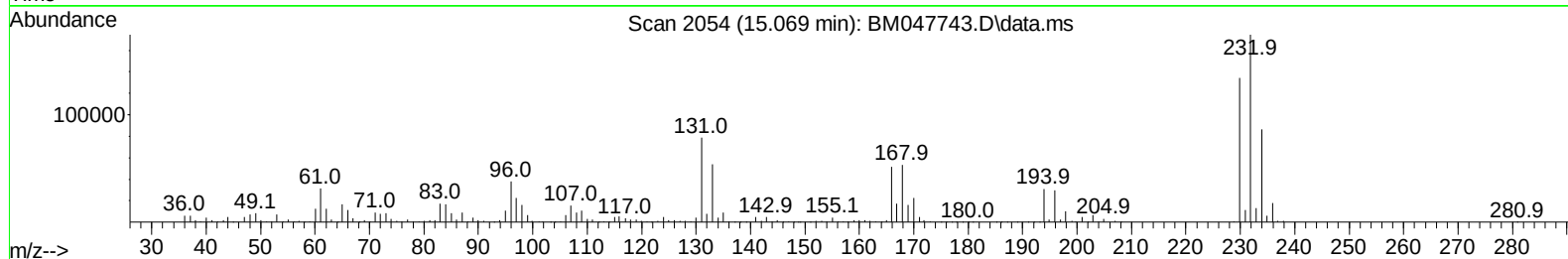
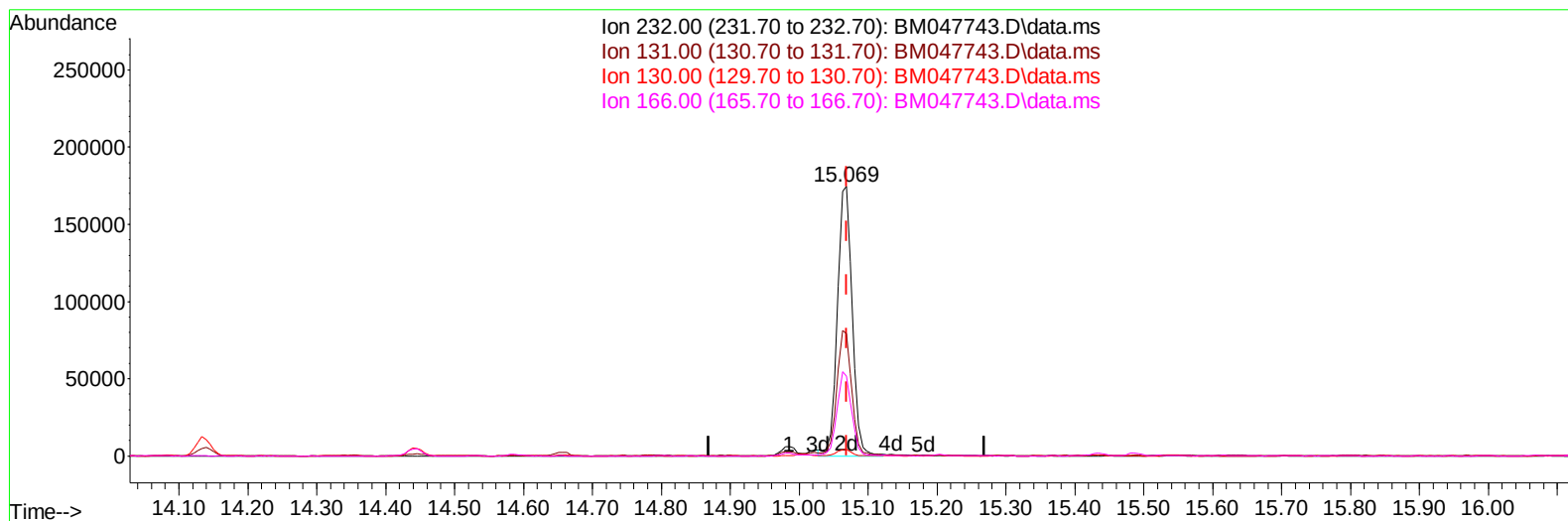
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TIC: BM047743.D\data.ms

(58) 2,3,4,6-Tetrachlorophenol

15.069min (+ 0.000) 19.29 ng/ul m

response 261236

Ion	Exp%	Act%
232.00	100.00	100.00
131.00	50.30	45.31
130.00	3.30	2.50#
166.00	31.60	29.59

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Quant Time: Oct 01 23:06:18 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.810	152	311506	20.000	ng/ul	0.00
20) Naphthalene-d8	10.604	136	1321328	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.445	164	834076	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.186	188	1603006	20.000	ng/ul	0.00
79) Chrysene-d12	21.403	240	1523533	20.000	ng/ul	0.00
88) Perylene-d12	24.409	264	1769785	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	23519	2.440	ng/uL	0.00
4) Pyridine-d5	3.675	84	392639	13.912	ng/ul	0.00
7) Phenol-d5	6.975	99	414663	14.077	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	260806	12.478	ng/ul	0.00
11) 2-Chlorophenol-d4	7.346	132	340151	15.872	ng/ul	0.00
15) 4-Methylphenol-d8	8.510	113	336268	15.354	ng/ul	0.00
21) Nitrobenzene-d5	8.963	128	165855	15.765	ng/ul	0.00
24) 2-Nitrophenol-d4	9.687	143	174353	17.294	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.222	165	329671	16.369	ng/ul	0.00
31) 4-Chloroaniline-d4	10.734	131	475345	15.267	ng/ul	0.00
46) Dimethylphthalate-d6	13.851	166	1102858	18.238	ng/ul	0.00
49) Acenaphthylene-d8	14.133	160	1274141	17.771	ng/ul	0.00
54) 4-Nitrophenol-d4	14.633	143	139827	11.614	ng/ul	0.00
60) Fluorene-d10	15.433	176	965268	18.429	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.545	200	111349	11.870	ng/ul	0.00
73) Anthracene-d10	17.280	188	1415224	17.955	ng/ul	0.00
81) Pyrene-d10	19.562	212	1698836	19.647	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.203	264	1748505	19.042	ng/ul	0.00
Target Compounds						
						Qvalue
58) 2,3,4,6-Tetrachlorophenol	15.069	232	261236m	19.292	ng/ul	
71) Pentachlorophenol	16.863	266	10931789	894.054	ng/ul	99

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

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