

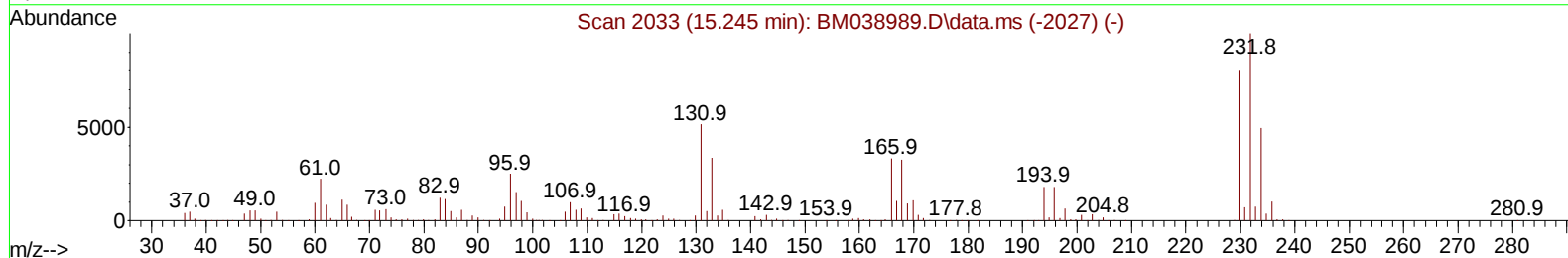
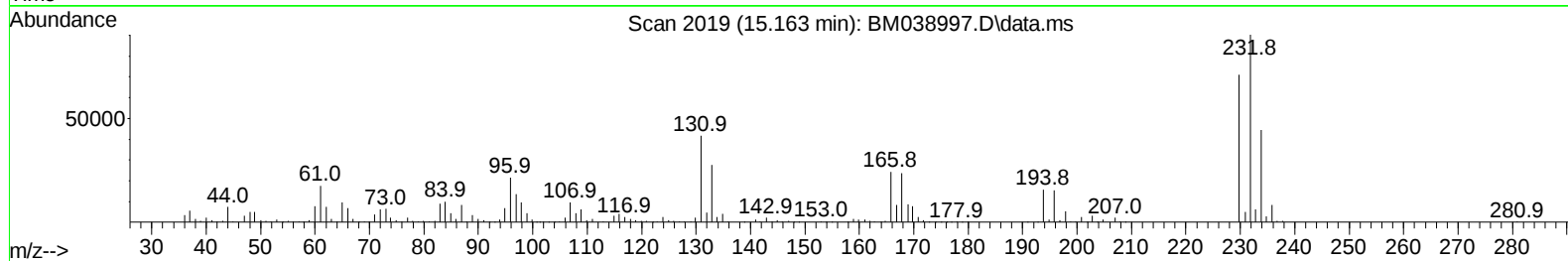
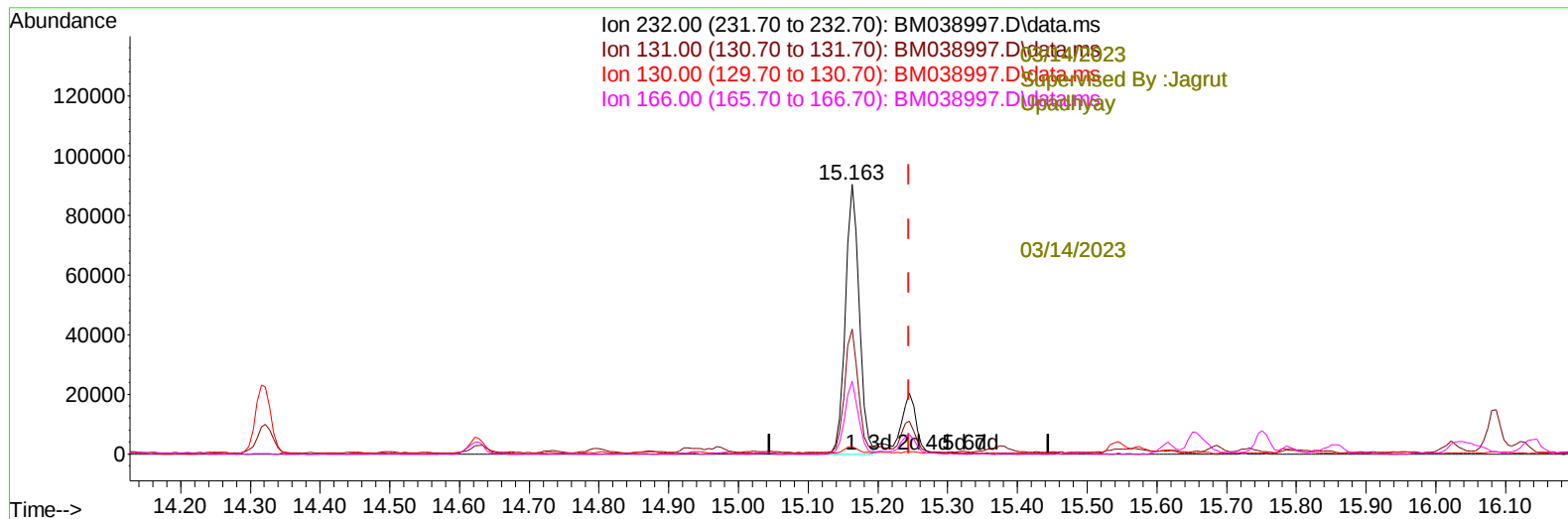
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM031323\
Data File : BM038997.D
Acq On : 13 Mar 2023 19:33
Operator : CG/JU
Sample : 01784-21
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DCAF6

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 03/14/2023
Supervised By :Jagrut Upadhyay 03/14/2023

Quant Time: Mar 14 01:43:29 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM031323.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Mar 13 16:25:27 2023
Response via : Initial Calibration



TIC: BM038997.D\data.ms

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

15.163min (-0.082) 8.21 ng/u1

response 123182

Ion	Exp%	Act%
232.00	100.00	100.00
131.00	51.60	46.33
130.00	2.80	2.71
166.00	33.10	27.01

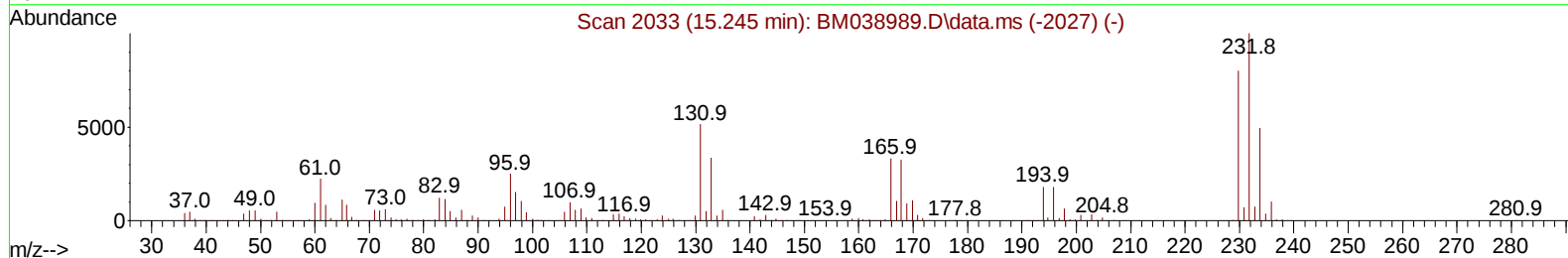
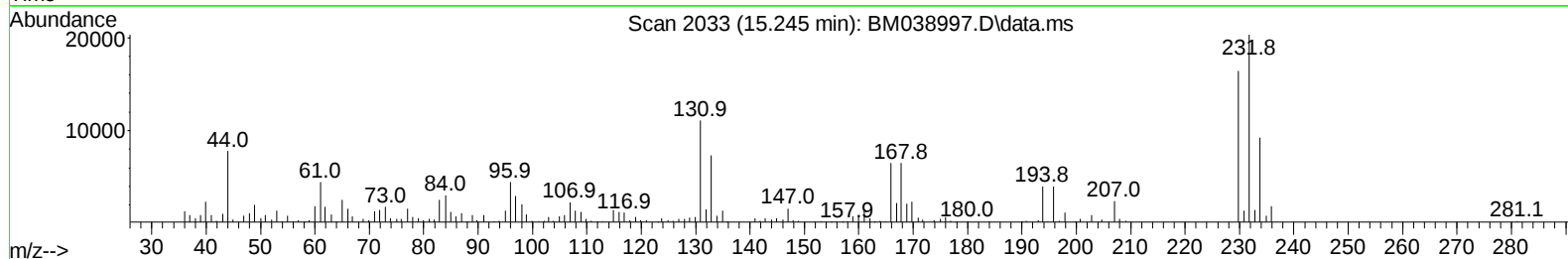
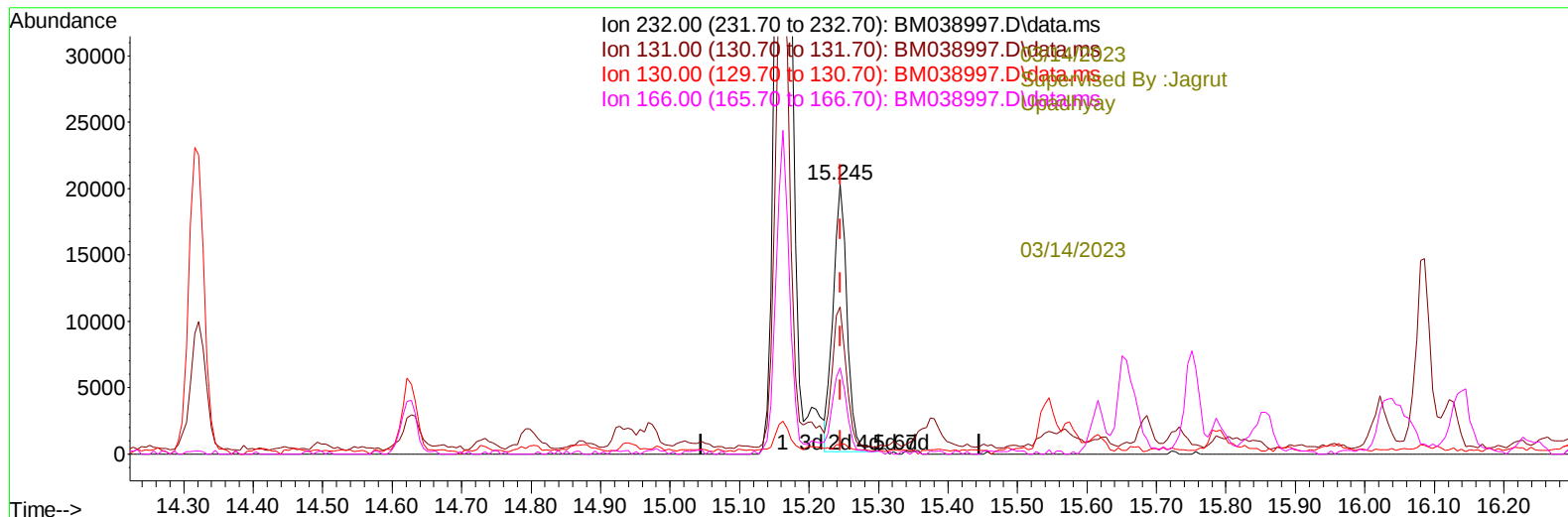
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TIC: BM038997.D\data.ms

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

15.245min (+ 0.000) 1.89 ng/ul m

response 28376

Ion	Exp%	Act%
232.00	100.00	100.00
131.00	51.60	54.56
130.00	2.80	3.46#
166.00	33.10	31.92

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)
-----03/14/2023							
Supervised By :Jagrut Upadhyay							
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.987	152		284595	20.000	ng/ul	0.00
20) Naphthalene-d8	10.798	136		1305159	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.627	164		844626	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.368	188		1837244	20.000	ng/ul	0.00
79) Chrysene-d12	21.550	240		1756734	20.000	ng/ul	0.00
88) Perylene-d12	23.997	264		1891571	20.000	ng/ul	0.00
-----03/14/2023							
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.369	96		34848	4.270	ng/uL	0.00
4) Pyridine-d5	3.799	84		164228	7.114	ng/ul	0.00
7) Phenol-d5	7.140	99		148248	5.411	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.310	67		520842	29.417	ng/ul	0.00
11) 2-Chlorophenol-d4	7.510	132		458439	22.479	ng/ul	0.00
15) 4-Methylphenol-d8	8.693	113		288805	13.475	ng/ul	0.00
21) Nitrobenzene-d5	9.151	128		323825	29.678	ng/ul	0.00
24) 2-Nitrophenol-d4	9.875	143		310061	28.204	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.416	165		525270	26.101	ng/ul	0.00
31) 4-Chloroaniline-d4	10.934	131		633497	19.463	ng/ul	0.00
46) Dimethylphthalate-d6	14.033	166		2082317	32.550	ng/ul	0.00
49) Acenaphthylene-d8	14.316	160		2198196	29.586	ng/ul	0.00
54) 4-Nitrophenol-d4	14.816	143		59437	4.459	ng/ul	0.00
60) Fluorene-d10	15.616	176		1779205	31.875	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.727	200		331477	28.486	ng/ul	0.00
73) Anthracene-d10	17.468	188		3093512	34.923	ng/ul	0.00
81) Pyrene-d10	19.756	212		3711510	36.086	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.839	264		3576718	36.833	ng/ul	0.00

Target Compounds							Qvalue
58) 2,3,4,6-Tetrachlorophenol	15.245	232		28376m	1.891	ng/ul	
71) Pentachlorophenol	17.015	266		140109	10.140	ng/ul	99

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

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