

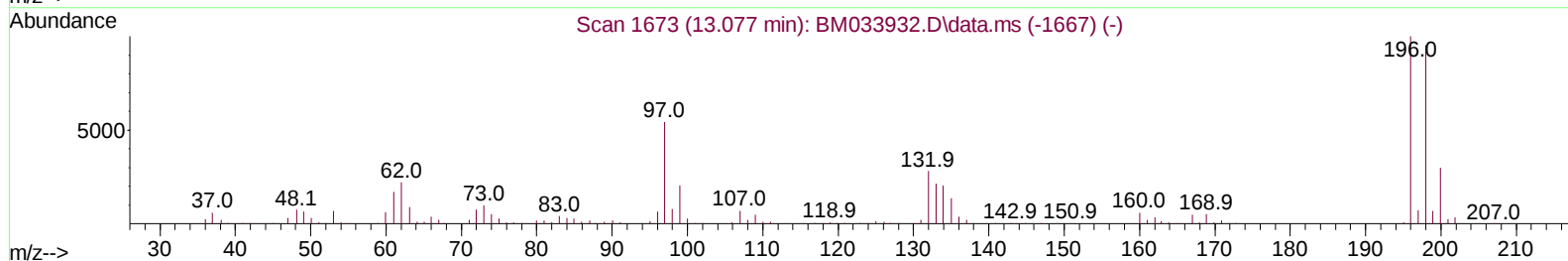
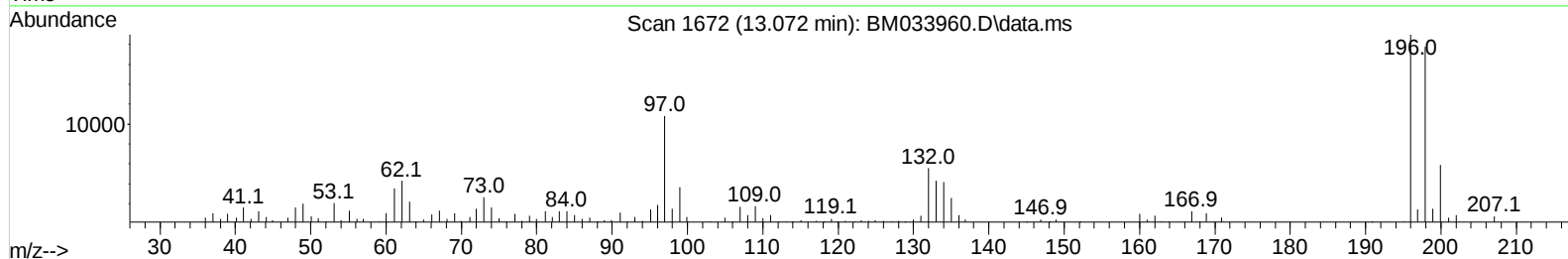
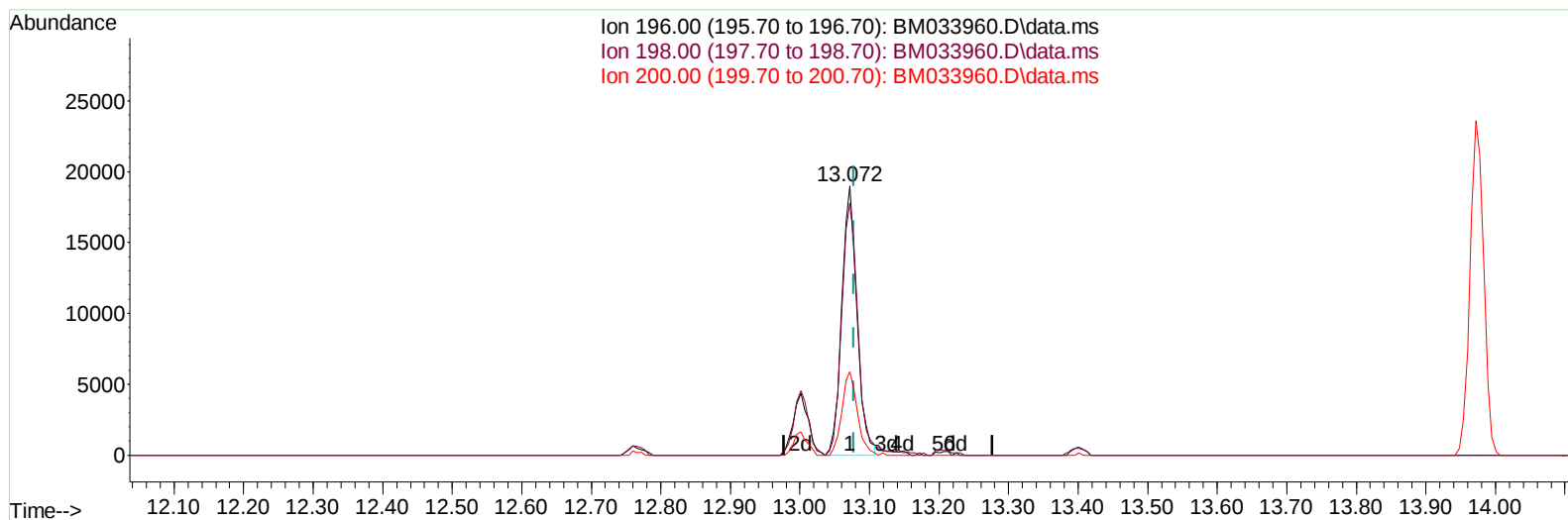
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
 Data File : BM033960.D
 Acq On : 02 Feb 2022 04:00
 Operator : CG/JU
 Sample : N1307-05
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0Q38

Manual IntegrationsAPPROVED

Quant Time: Feb 02 05:59:42 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Feb 01 02:16:28 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/02/2022
 Supervised By :Yogesh Patel 02/04/2022



TIC: BM033960.D\data.ms

(42) 2,4,5-Trichlorophenol

13.072min (-0.006) 4.39 ng/u1

response 29230

Ion	Exp%	Act%
196.00	100.00	100.00
198.00	102.70	93.64
200.00	34.90	30.94
0.00	0.00	0.00

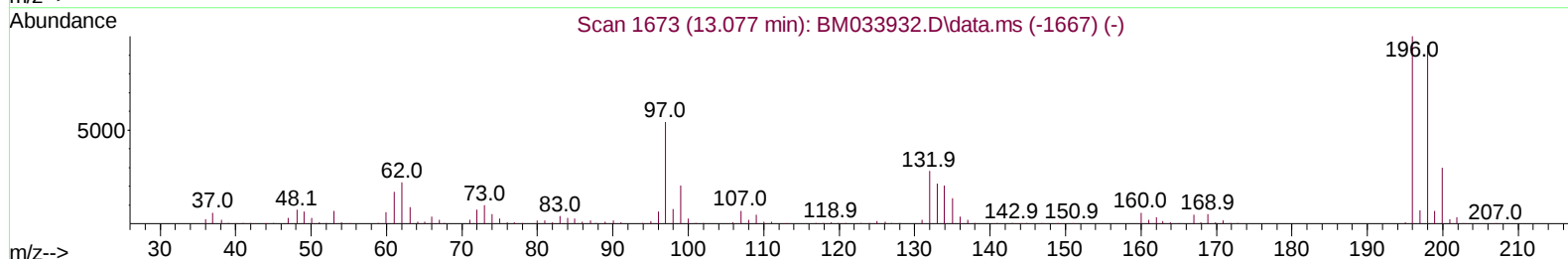
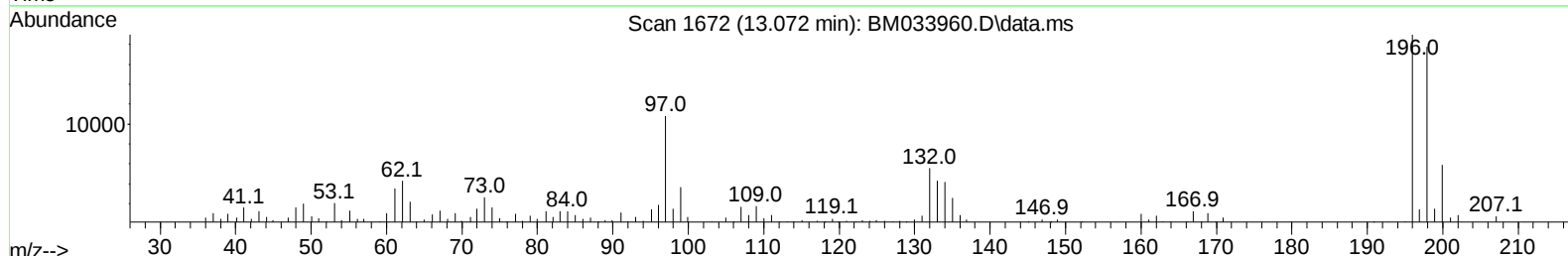
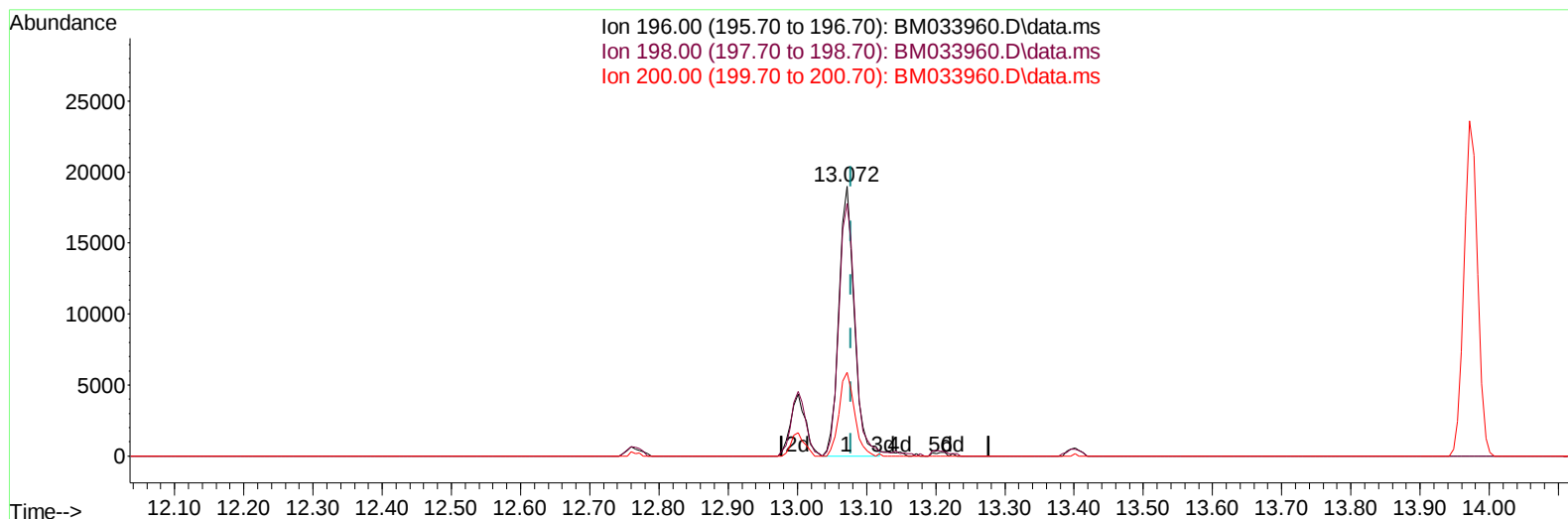
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TIC: BM033960.D\data.ms

(42) 2,4,5-Trichlorophenol

13.072min (-0.006) 4.45 ng/ul m

response 29571

Ion	Exp%	Act%
196.00	100.00	100.00
198.00	102.70	93.64
200.00	34.90	30.94
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.925	152	128114	20.000	ng/ul	0.00
20) Naphthalene-d8	10.737	136	530474	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.566	164	316646	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.307	188	590339	20.000	ng/ul	0.00
79) Chrysene-d12	21.459	240	510690	20.000	ng/ul	0.00
88) Perylene-d12	23.847	264	513929	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.290	96	13071	4.024	ng/uL	0.00
4) Pyridine-d5	3.719	84	100744	11.111	ng/ul	0.00
7) Phenol-d5	7.084	99	68997	6.429	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.248	67	211147	32.269	ng/ul	0.00
11) 2-Chlorophenol-d4	7.454	132	216358	25.161	ng/ul	0.00
15) 4-Methylphenol-d8	8.631	113	135787	15.966	ng/ul	-0.01
21) Nitrobenzene-d5	9.084	128	136883	32.159	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.813	143	146826	31.943	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.354	165	230905	27.760	ng/ul	0.00
31) 4-Chloroaniline-d4	10.866	131	142591	12.414	ng/ul	0.00
46) Dimethylphthalate-d6	13.972	166	813528	34.381	ng/ul	0.00
49) Acenaphthylene-d8	14.254	160	985833	34.010	ng/ul	0.00
54) 4-Nitrophenol-d4	14.754	143	20245	5.008	ng/ul	0.00
60) Fluorene-d10	15.548	176	666597	33.739	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.666	200	147371	39.764	ng/ul	0.00
73) Anthracene-d10	17.401	188	1020191	36.086	ng/ul	0.00
81) Pyrene-d10	19.683	212	1063704	35.280	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.694	264	969316	35.909	ng/ul	-0.01
Target Compounds						
					Qvalue	
41) 2,4,6-Trichlorophenol	13.001	196	6506	1.037	ng/ul#	92
42) 2,4,5-Trichlorophenol	13.072	196	29571m	4.445	ng/ul	
58) 2,3,4,6-Tetrachlorophenol	15.183	232	232356	42.453	ng/ul#	98
71) Pentachlorophenol	16.983	266	8933073	1909.662	ng/ul	94

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

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