

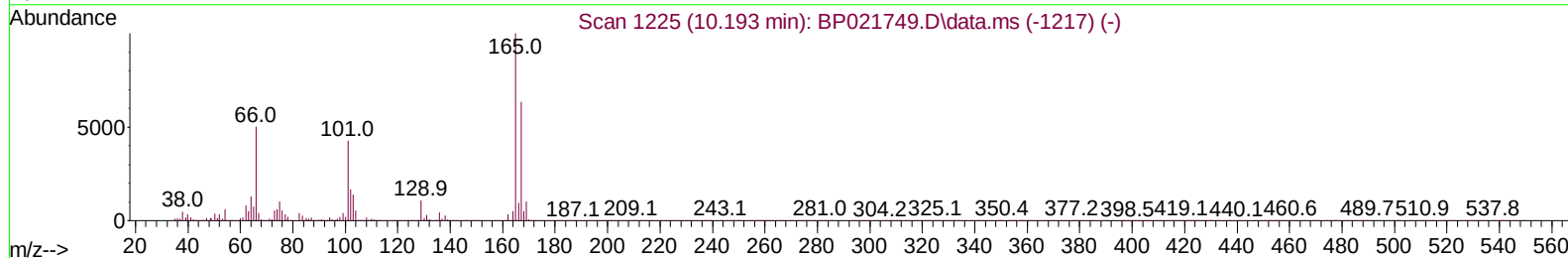
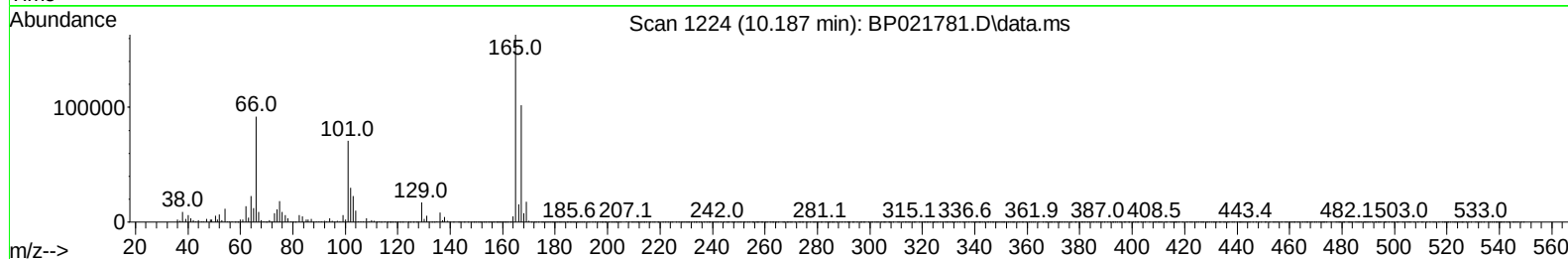
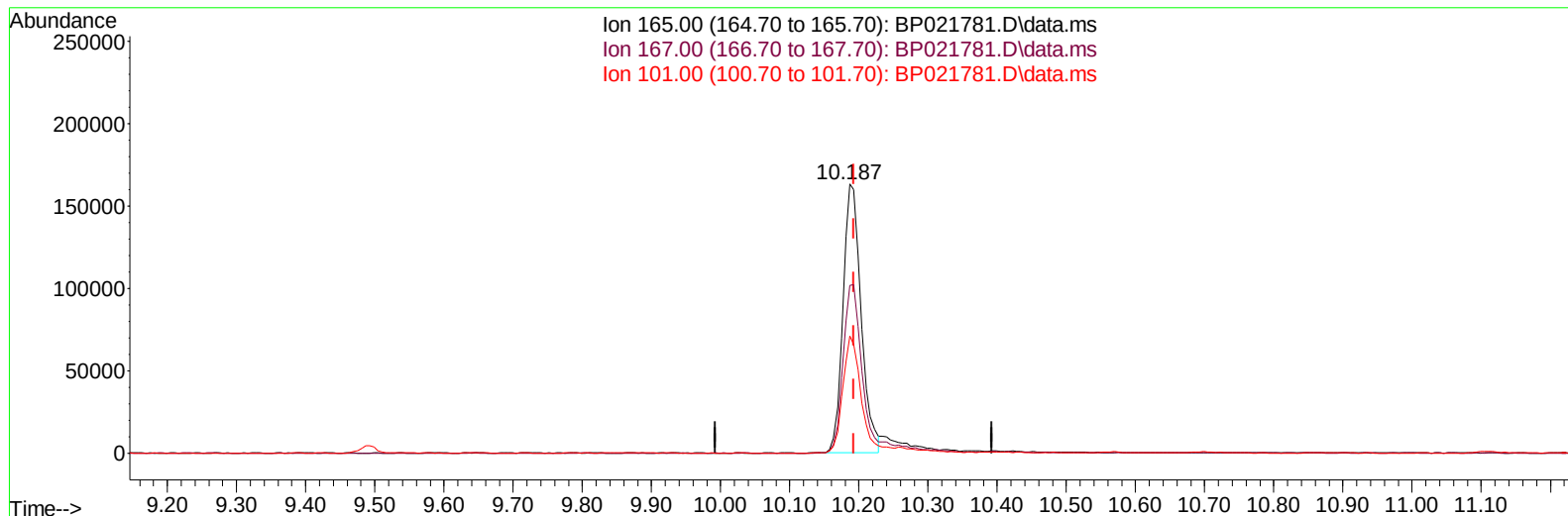
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP083024\
Data File : BP021781.D
Acq On : 31 Aug 2024 11:52
Operator : RC/JU
Sample : P3733-17
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44A9

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 09/03/2024
Supervised By :mohammad ahmed 09/03/2024

Quant Time: Sep 02 01:09:47 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Aug 29 23:56:59 2024
Response via : Initial Calibration



TIC: BP021781.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.187min (-0.006) 19.82 ng/ul

response 298115

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.10	62.37
101.00	41.40	43.55
0.00	0.00	0.00

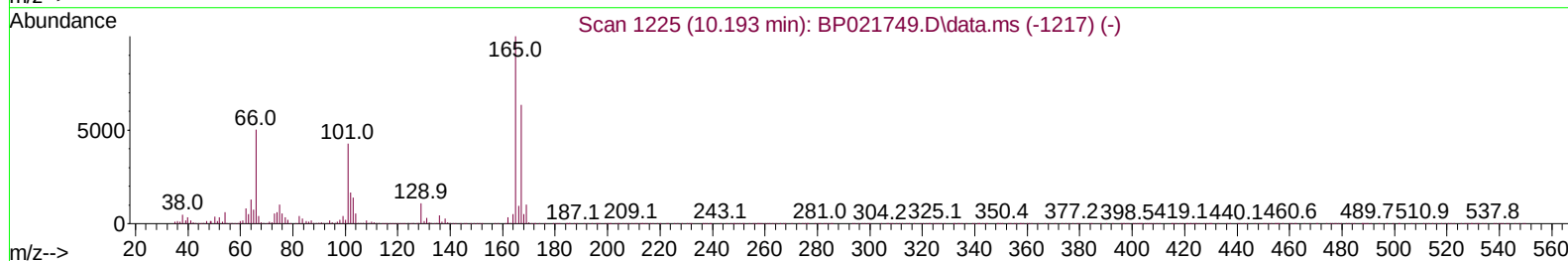
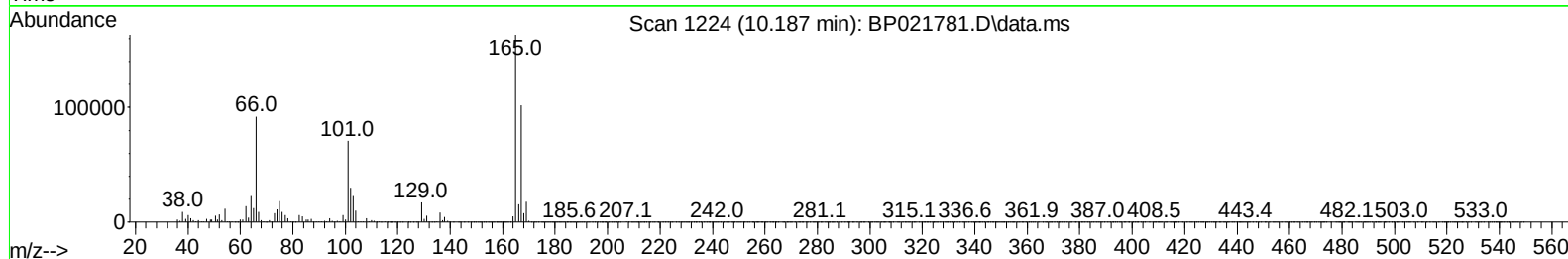
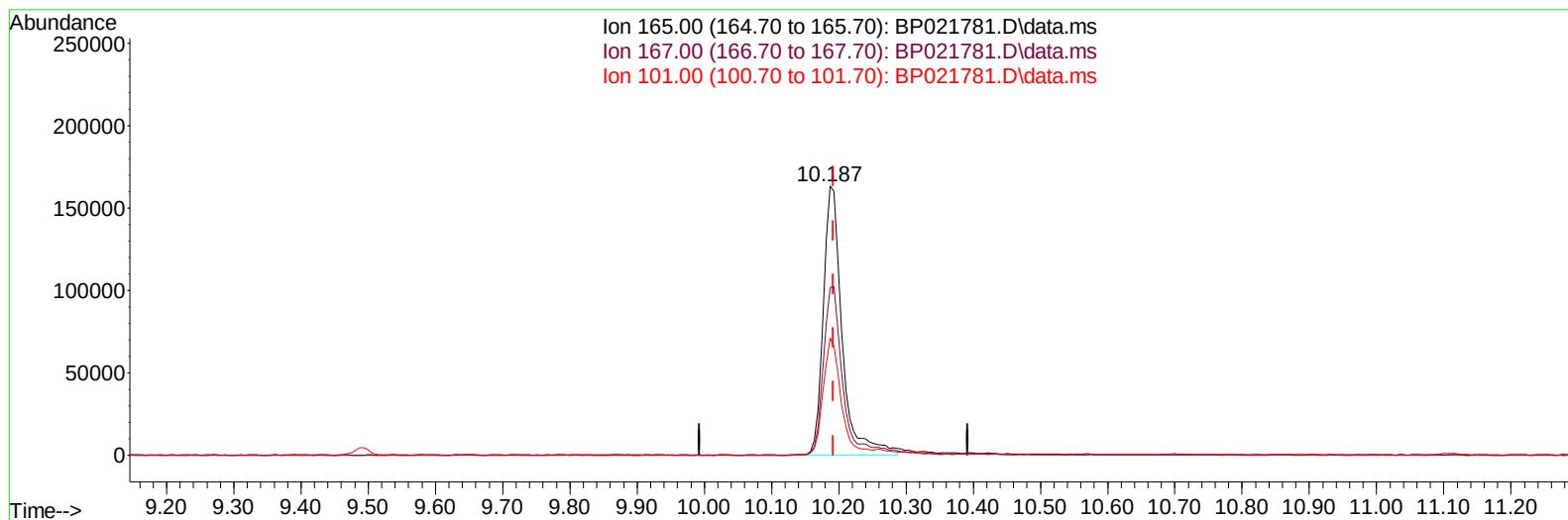
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TIC: BP021781.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.187min (-0.006) 21.41 ng/ul m

response 322121

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.10	62.37
101.00	41.40	43.55
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 09/03/2024
 Supervised By :mohammad ahmed 09/03/2024

Quant Time: Sep 02 01:55:04 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 29 23:56:59 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	248586	20.000	ng/ul	0.00
20) Naphthalene-d8	10.569	136	934813	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.416	164	555591	20.000	ng/ul	-0.02
64) Phenanthrene-d10	17.222	188	1151840	20.000	ng/ul	0.00
79) Chrysene-d12	21.680	240	1255935	20.000	ng/ul	0.01
88) Perylene-d12	25.068	264	1541802	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	31452	4.200	ng/uL	0.00
4) Pyridine-d5	3.687	84	431972	19.946	ng/ul	0.00
7) Phenol-d5	6.952	99	565406	21.313	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.111	67	313030	21.682	ng/ul	0.00
11) 2-Chlorophenol-d4	7.322	132	377498	22.041	ng/ul	0.00
15) 4-Methylphenol-d8	8.481	113	424020	20.684	ng/ul	0.00
21) Nitrobenzene-d5	8.928	128	174541	23.207	ng/ul	0.00
24) 2-Nitrophenol-d4	9.646	143	172120	22.546	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.187	165	322121m	21.412	ng/ul	0.00
31) 4-Chloroaniline-d4	10.699	131	491188	22.442	ng/ul	0.00
46) Dimethylphthalate-d6	13.822	166	1046837	24.624	ng/ul	-0.02
49) Acenaphthylene-d8	14.110	160	1192112	24.976	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.634	143	162714	20.118	ng/ul	0.00
60) Fluorene-d10	15.428	176	914670	25.592	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.545	200	105911	16.718	ng/ul	-0.01
73) Anthracene-d10	17.322	188	1404575	25.694	ng/ul	-0.02
81) Pyrene-d10	19.692	212	1794149	24.990	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.839	264	2158959	26.260	ng/ul	0.00

Target Compounds Qvalue

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

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