

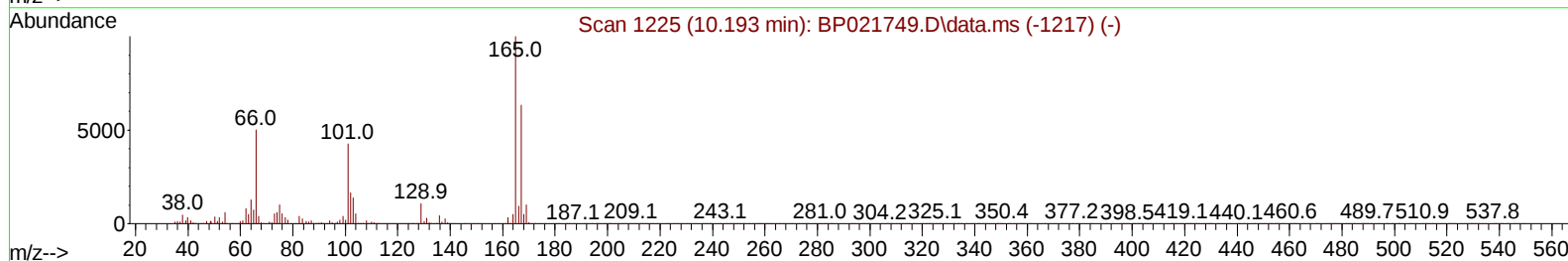
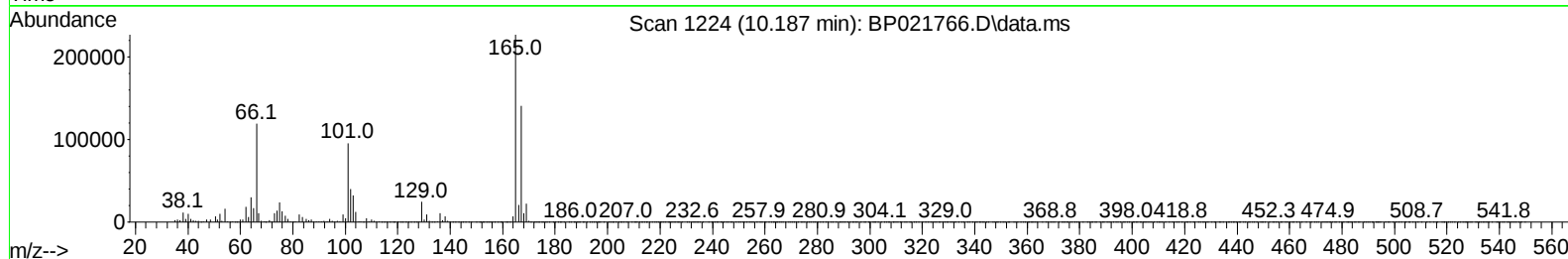
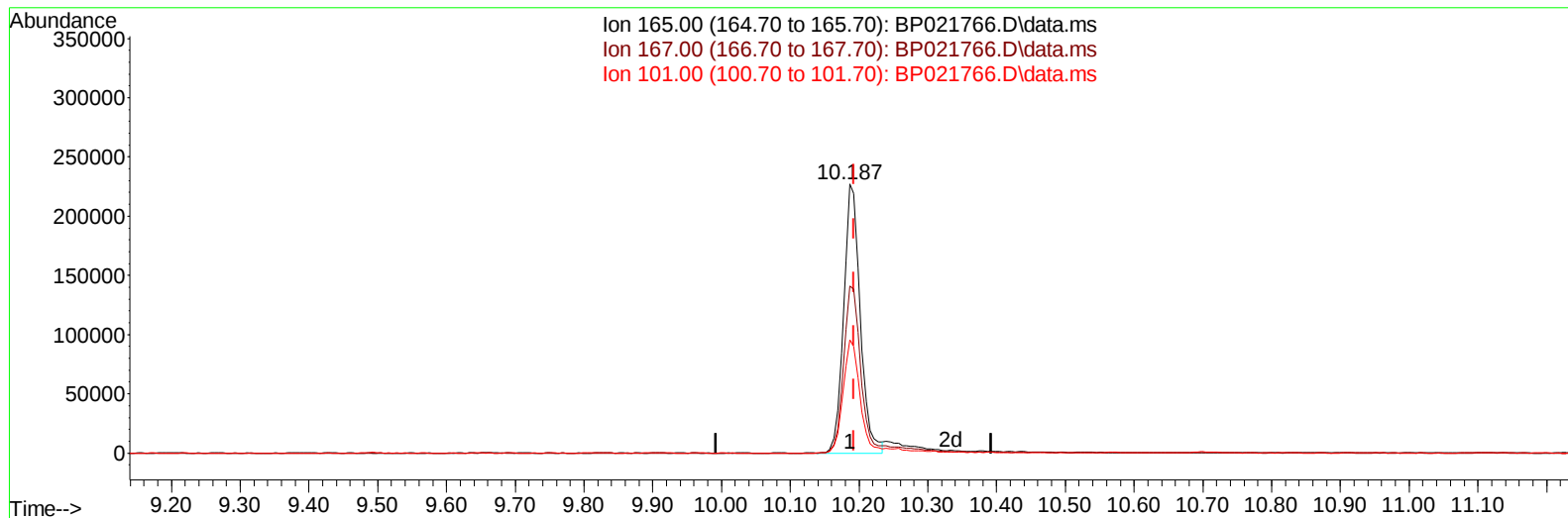
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP083024\
Data File : BP021766.D
Acq On : 31 Aug 2024 00:15
Operator : RC/JU
Sample : P3733-06
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44A0

Manual IntegrationsAPPROVED

Quant Time: Aug 31 00:44:54 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

Reviewed By :Yogesh Patel 09/02/2024
Supervised By :mohammad ahmed 09/03/2024



TIC: BP021766.D\data.ms

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

10.187min (-0.006) 22.73 ng/u1

response 381561

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.10	62.08
101.00	41.40	42.10
0.00	0.00	0.00

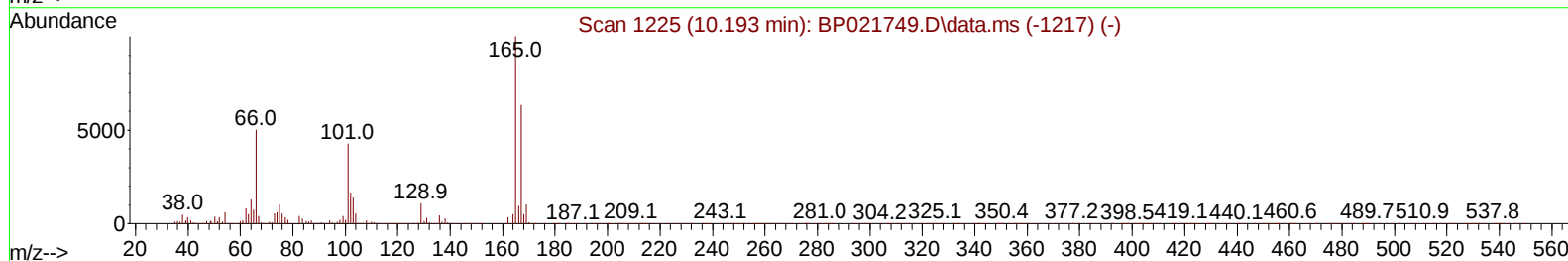
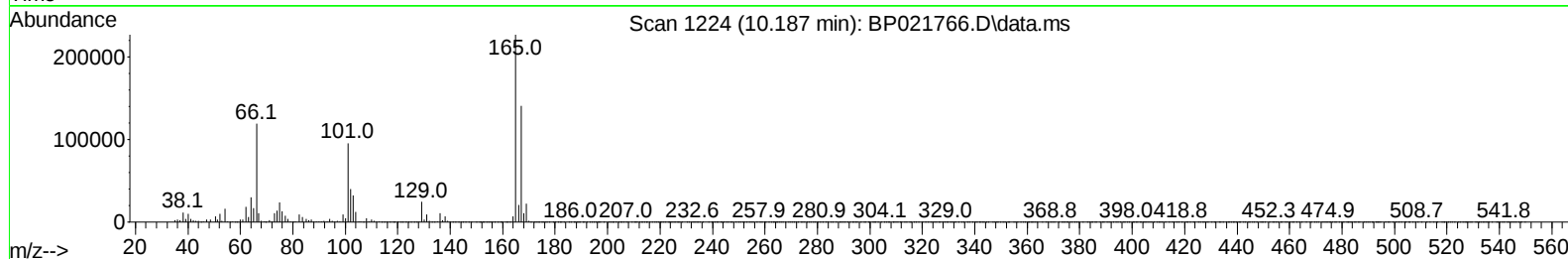
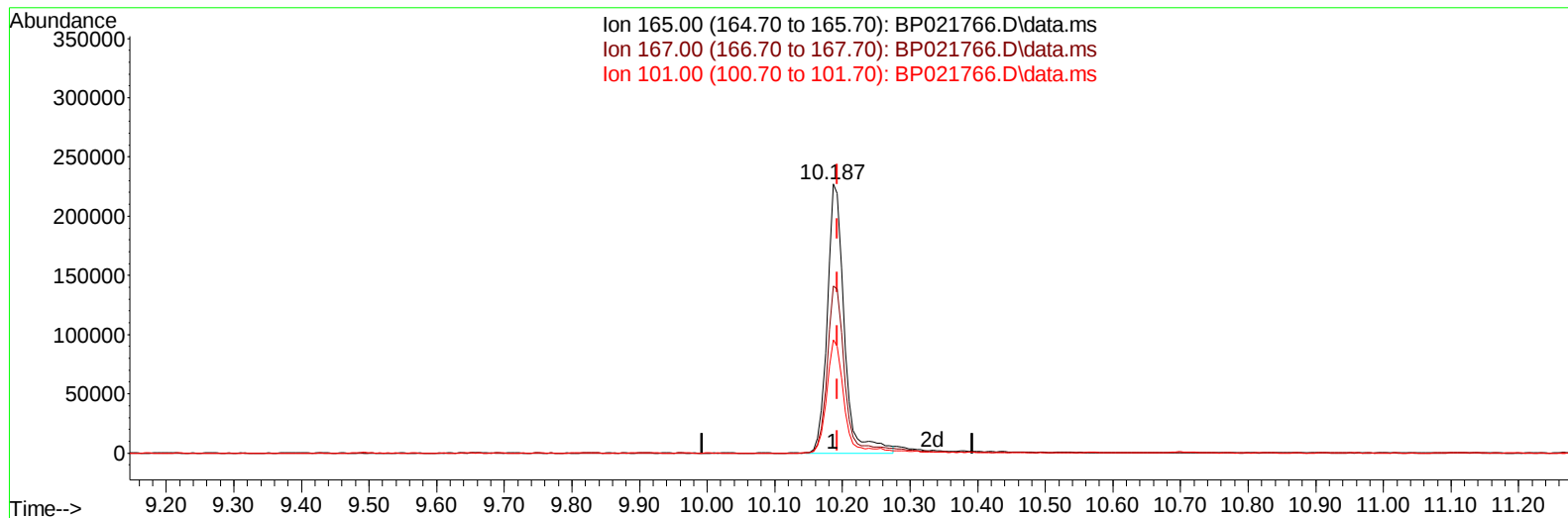
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TIC: BP021766.D\data.ms

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

10.187min (-0.006) 23.86 ng/ul m

response 400535

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.10	62.08
101.00	41.40	42.10
0.00	0.00	0.00

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 Sample : P3733-06
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 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A44A0

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 09/02/2024
 Supervised By :mohammad ahmed 09/03/2024

Quant Time: Aug 31 00:47:31 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	272849	20.000	ng/ul	0.00
20) Naphthalene-d8	10.569	136	1042956	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.428	164	590400	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.227	188	1086115	20.000	ng/ul	0.00
79) Chrysene-d12	21.668	240	1105790	20.000	ng/ul	0.00
88) Perylene-d12	25.062	264	1421901	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	37289	4.537	ng/uL	0.00
4) Pyridine-d5	3.687	84	537770	22.623	ng/ul	0.00
7) Phenol-d5	6.957	99	705069	24.214	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.110	67	380536	24.014	ng/ul	0.00
11) 2-Chlorophenol-d4	7.322	132	469067	24.951	ng/ul	0.00
15) 4-Methylphenol-d8	8.481	113	527838	23.459	ng/ul	0.00
21) Nitrobenzene-d5	8.928	128	220978	26.335	ng/ul	0.00
24) 2-Nitrophenol-d4	9.651	143	225829	26.514	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.187	165	400535m	23.863	ng/ul	0.00
31) 4-Chloroaniline-d4	10.698	131	574415	23.523	ng/ul	0.00
46) Dimethylphthalate-d6	13.828	166	1205569	26.686	ng/ul	-0.01
49) Acenaphthylene-d8	14.116	160	1424357	28.083	ng/ul	0.00
54) 4-Nitrophenol-d4	14.628	143	151704	17.651	ng/ul	0.00
60) Fluorene-d10	15.433	176	1020877	26.879	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.545	200	103153	17.268	ng/ul	-0.01
73) Anthracene-d10	17.333	188	1482594	28.763	ng/ul	0.00
81) Pyrene-d10	19.692	212	1749110	27.671	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.827	264	2213779	29.198	ng/ul	-0.01

Target Compounds Qvalue

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

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