

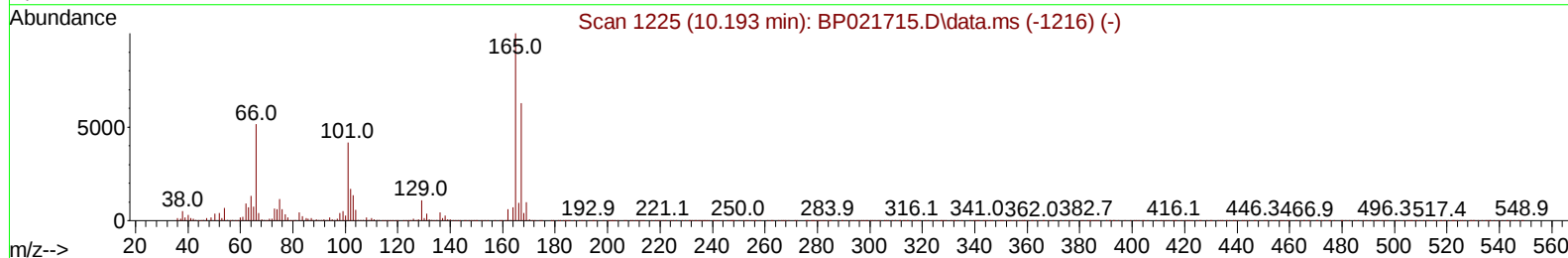
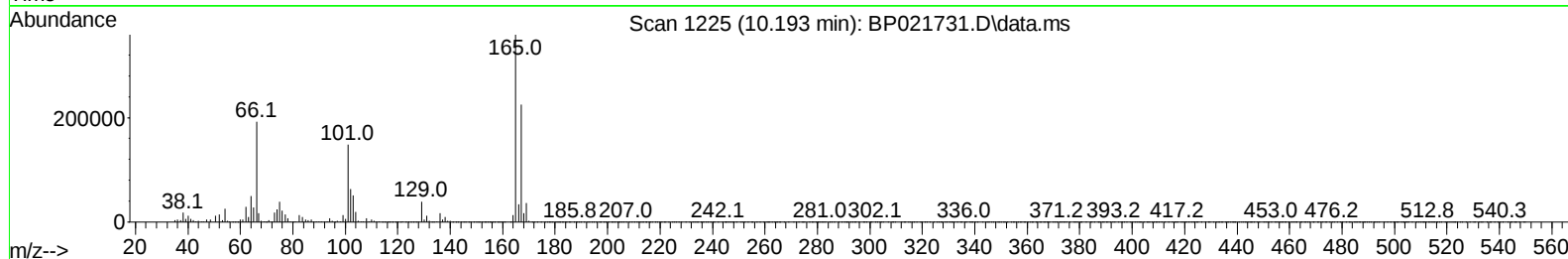
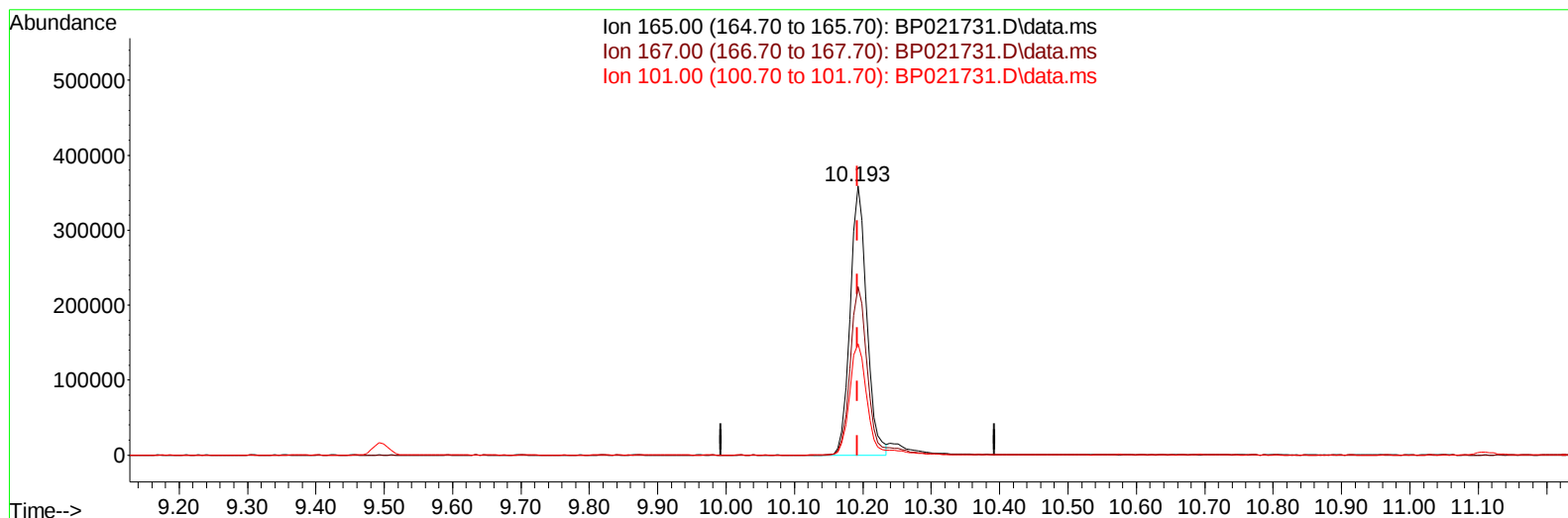
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082924\
Data File : BP021731.D
Acq On : 29 Aug 2024 21:51
Operator : RC/JU
Sample : P3721-06
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCYJ0

Manual IntegrationsAPPROVED

Quant Time: Aug 29 23:59:32 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 :Jagrut Upadhyay 08/30/2024
Supervised By :mohammad ahmed 09/02/2024



TIC: BP021731.D\data.ms

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

10.193min (-0.000) 26.53 ng/u1

response 605115

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.10	62.59
101.00	41.40	41.47
0.00	0.00	0.00

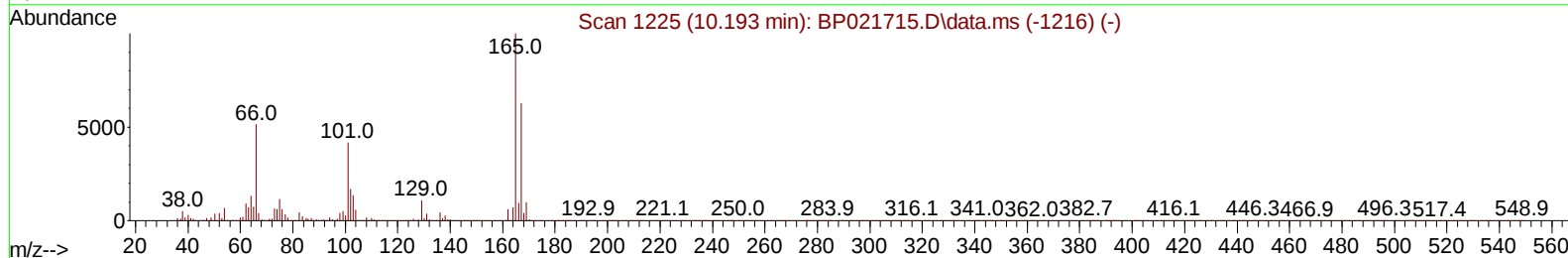
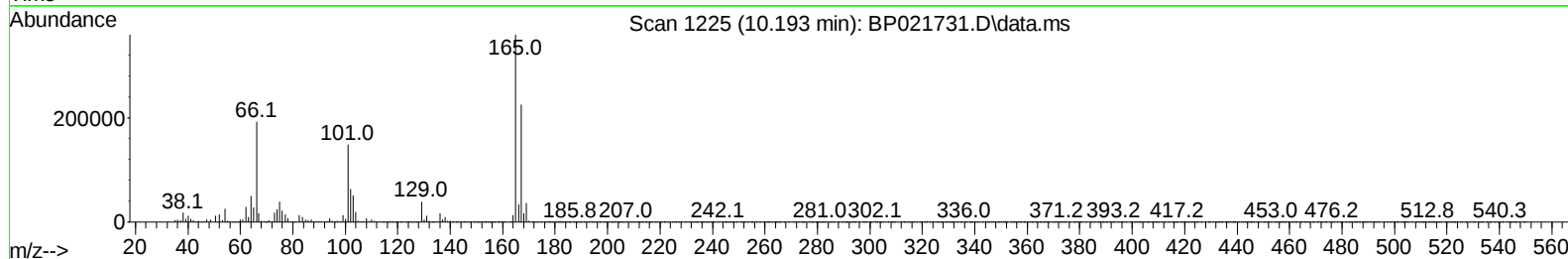
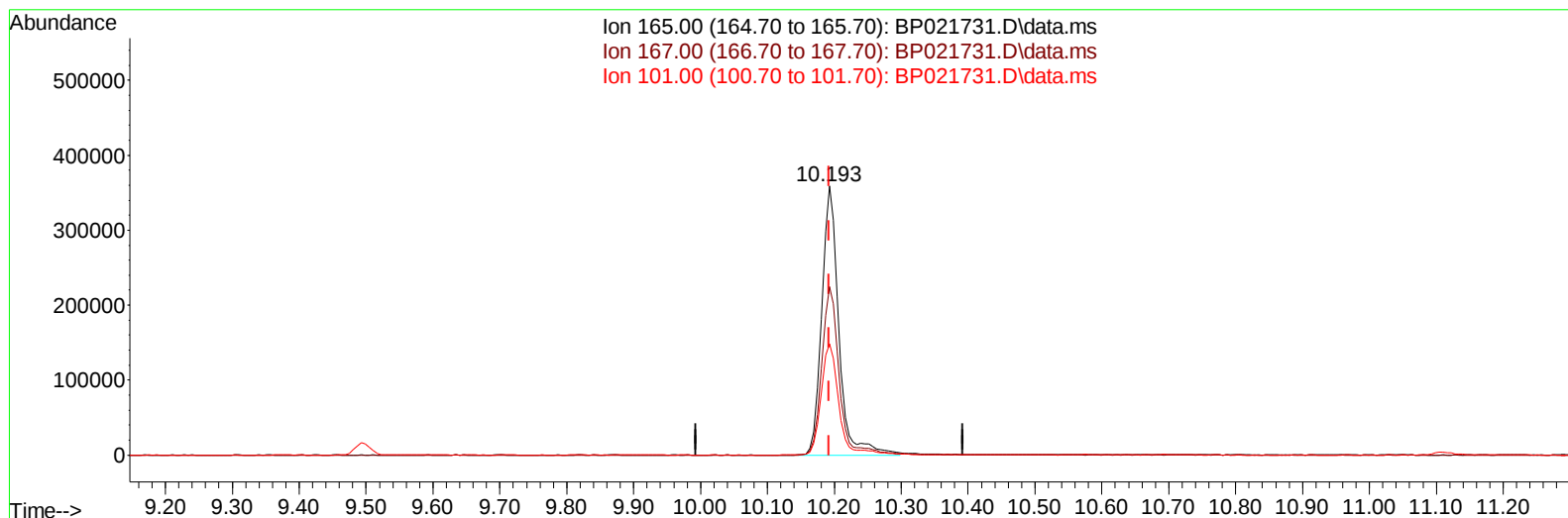
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TIC: BP021731.D\data.ms

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

10.193min (-0.000) 28.00 ng/ul m

response 638775

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.10	62.59
101.00	41.40	41.47
0.00	0.00	0.00

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 BNA_P
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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/30/2024
 Supervised By :mohammad ahmed 09/02/2024

Quant Time: Aug 30 00:04:28 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	7.787	152	334532	20.000	ng/ul	0.00
20) Naphthalene-d8	10.569	136	1417444	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.434	164	940347	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.233	188	2057270	20.000	ng/ul	0.00
79) Chrysene-d12	21.674	240	2088248	20.000	ng/ul	0.00
88) Perylene-d12	25.074	264	2388469	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	43172	4.284	ng/uL	0.00
4) Pyridine-d5	3.687	84	177954	6.106	ng/ul	0.00
7) Phenol-d5	6.957	99	977136	27.370	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.116	67	516464	26.582	ng/ul	0.00
11) 2-Chlorophenol-d4	7.322	132	631868	27.414	ng/ul	0.00
15) 4-Methylphenol-d8	8.487	113	759884	27.544	ng/ul	0.00
21) Nitrobenzene-d5	8.934	128	308714	27.071	ng/ul	0.00
24) 2-Nitrophenol-d4	9.651	143	316405	27.334	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.193	165	638775m	28.003	ng/ul	0.00
31) 4-Chloroaniline-d4	10.698	131	428172	12.902	ng/ul	0.00
46) Dimethylphthalate-d6	13.828	166	1977529	27.484	ng/ul	-0.01
49) Acenaphthylene-d8	14.116	160	2275198	28.164	ng/ul	0.00
54) 4-Nitrophenol-d4	14.628	143	335186	24.486	ng/ul	0.00
60) Fluorene-d10	15.439	176	1729617	28.592	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.551	200	240471	21.252	ng/ul	0.00
73) Anthracene-d10	17.339	188	2799988	28.678	ng/ul	0.00
81) Pyrene-d10	19.704	212	3380312	28.317	ng/ul	0.01
92) Benzo(a)pyrene-d12	24.845	264	3626803	28.476	ng/ul	0.00

Target Compounds Qvalue

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

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