

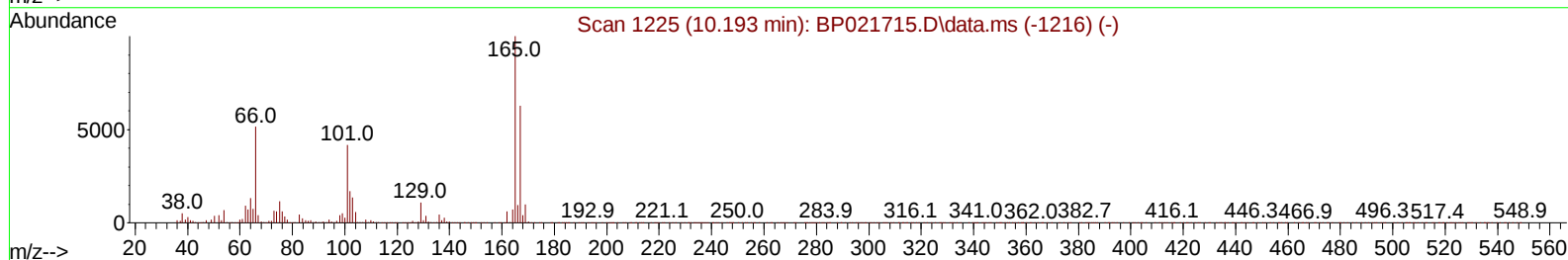
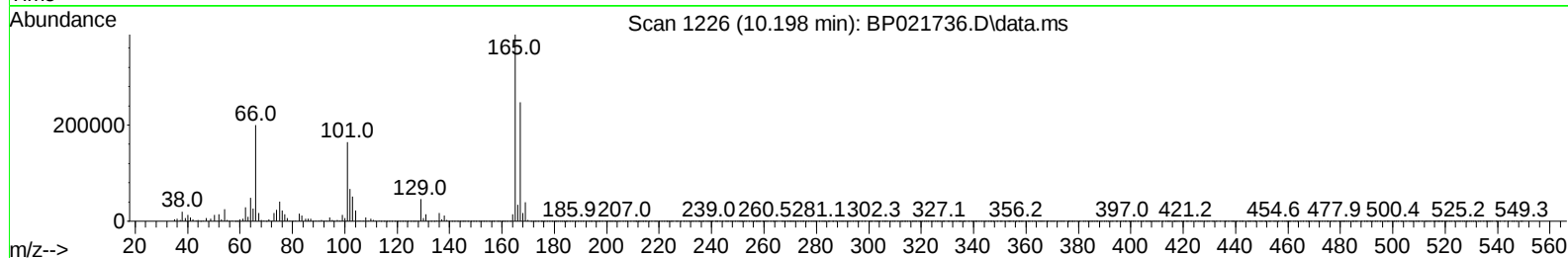
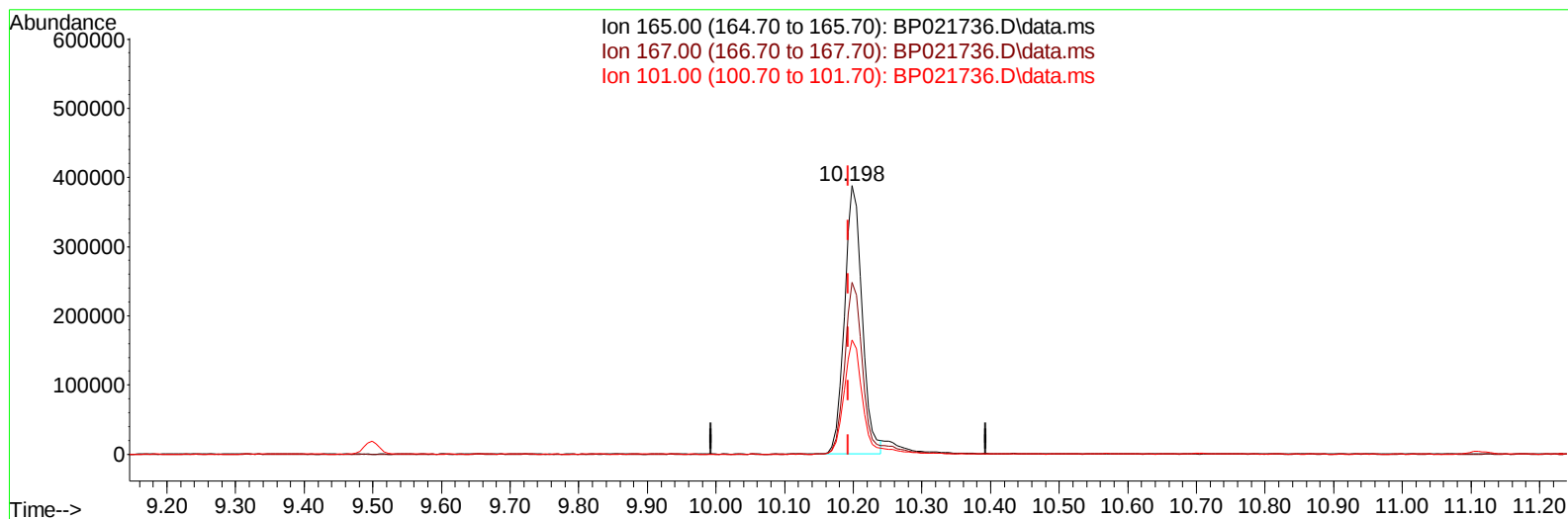
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082924\
Data File : BP021736.D
Acq On : 30 Aug 2024 01:27
Operator : RC/JU
Sample : P3721-07
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCXZ0

Manual IntegrationsAPPROVED

Quant Time: Aug 30 01:53:09 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: BP021736.D\data.ms

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

10.198min (+ 0.006) 25.49 ng/u1

response 688662

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.10	63.91
101.00	41.40	42.49
0.00	0.00	0.00

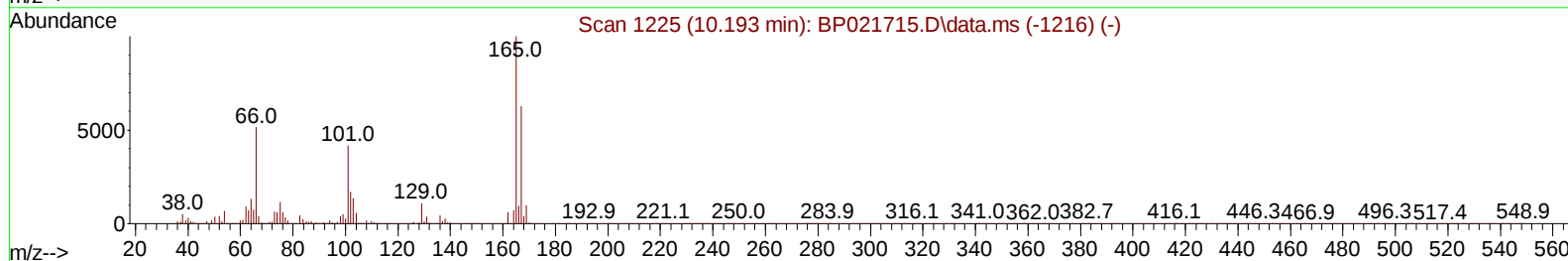
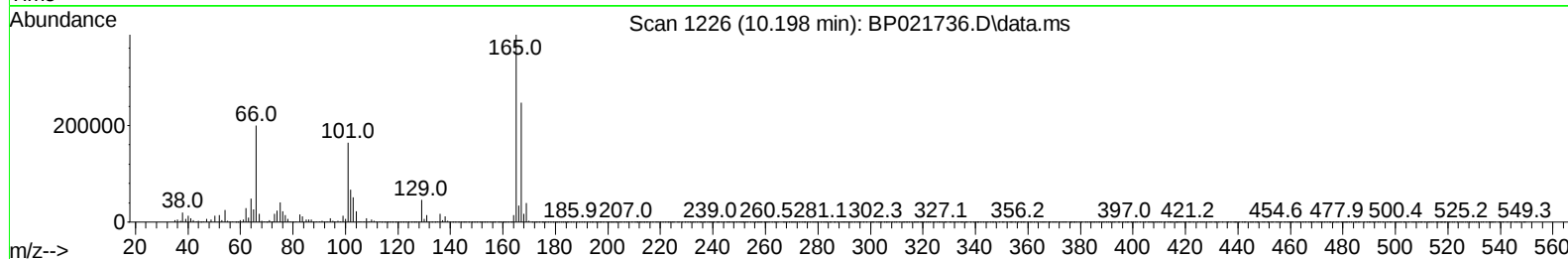
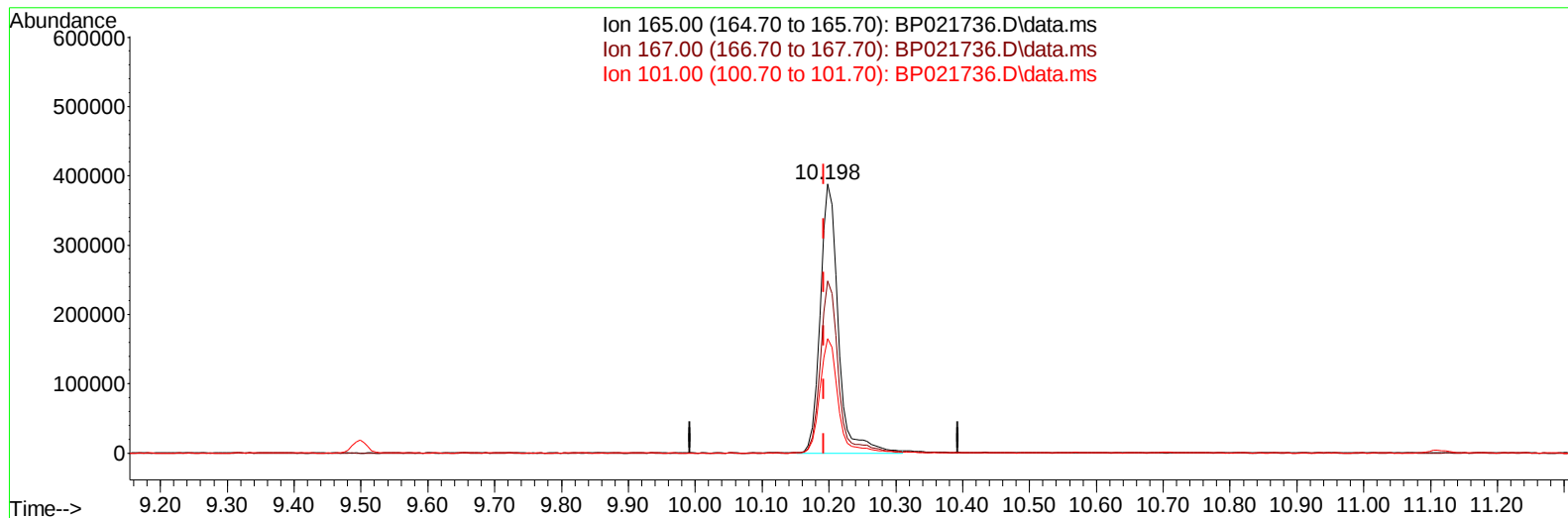
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TIC: BP021736.D\data.ms

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

10.198min (+ 0.006) 27.01 ng/ul m

response 729574

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.10	63.91
101.00	41.40	42.49
0.00	0.00	0.00

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 QLast Update : Thu Aug 29 23:56:59 2024
 Response via : Initial Calibration

Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.787	152		384817	20.000	ng/ul	0.00
20) Naphthalene-d8	10.575	136		1678502	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.428	164		1130860	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.233	188		2402309	20.000	ng/ul	0.00
79) Chrysene-d12	21.680	240		2115873	20.000	ng/ul	0.01
88) Perylene-d12	25.086	264		2378032	20.000	ng/ul	0.02
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.275	96		46171	3.983	ng/uL	0.00
4) Pyridine-d5	3.693	84		293105	8.743	ng/ul	0.00
7) Phenol-d5	6.963	99		1090177	26.546	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.116	67		568974	25.458	ng/ul	0.00
11) 2-Chlorophenol-d4	7.328	132		717690	27.069	ng/ul	0.00
15) 4-Methylphenol-d8	8.493	113		827931	26.089	ng/ul	0.00
21) Nitrobenzene-d5	8.934	128		358983	26.583	ng/ul	0.00
24) 2-Nitrophenol-d4	9.657	143		392049	28.601	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.198	165		729574m	27.009	ng/ul	0.00
31) 4-Chloroaniline-d4	10.704	131		646117	16.441	ng/ul	0.00
46) Dimethylphthalate-d6	13.834	166		2307310	26.665	ng/ul	0.00
49) Acenaphthylene-d8	14.122	160		2630691	27.079	ng/ul	0.00
54) 4-Nitrophenol-d4	14.639	143		400095	24.303	ng/ul	0.01
60) Fluorene-d10	15.439	176		1989004	27.341	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200		274957	20.810	ng/ul	0.00
73) Anthracene-d10	17.333	188		2929357	25.694	ng/ul	0.00
81) Pyrene-d10	19.704	212		3553934	29.383	ng/ul	0.01
92) Benzo(a)pyrene-d12	24.851	264		3438217	27.114	ng/ul	0.01

Target Compounds Qvalue

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

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