Quantitation Report (Qedit)

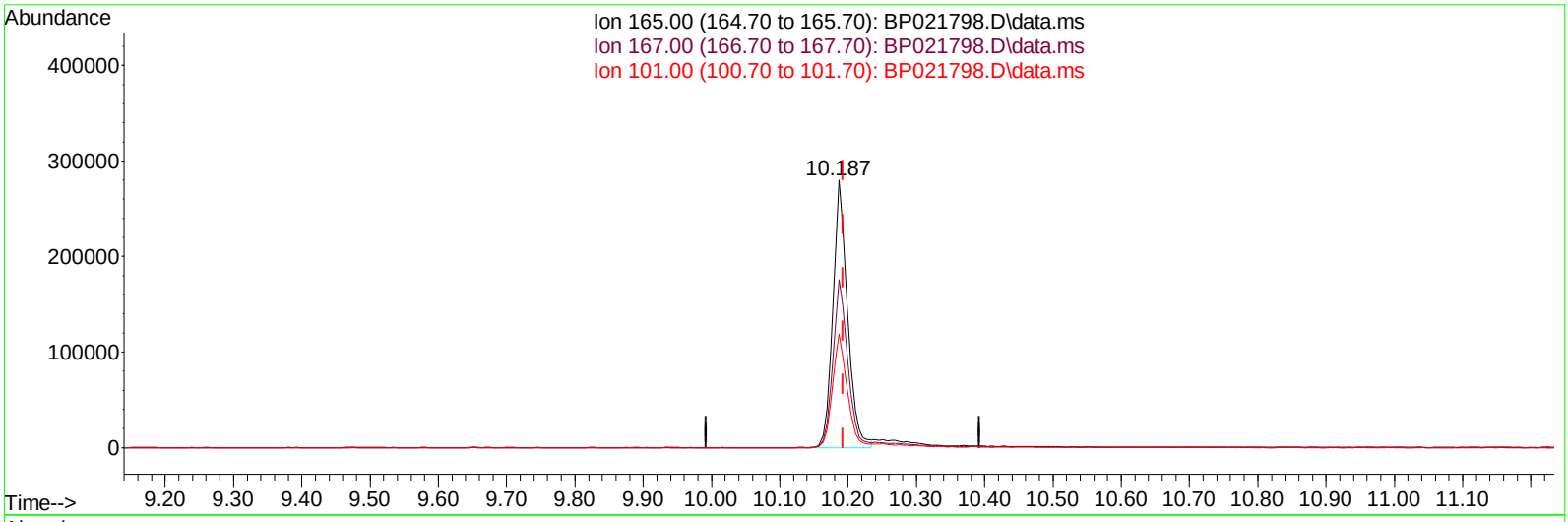
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP090324\
Data File : BP021798.D
Acq On : 03 Sep 2024 22:05
Operator : RC/JU
Sample : P3785-03
Misc :
ALS Vial : 14 Sample Multiplier: 1

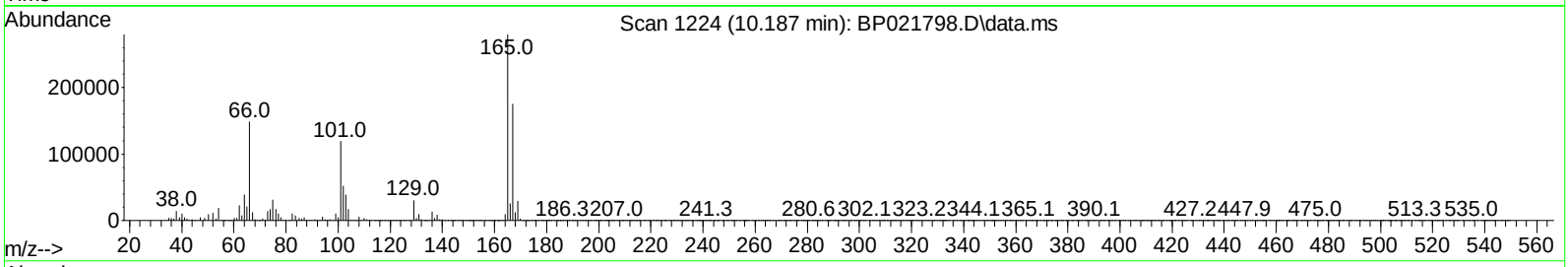
Quant Time: Sep 04 04:03:34 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Aug 29 23:56:59 2024
Response via : Initial Calibration

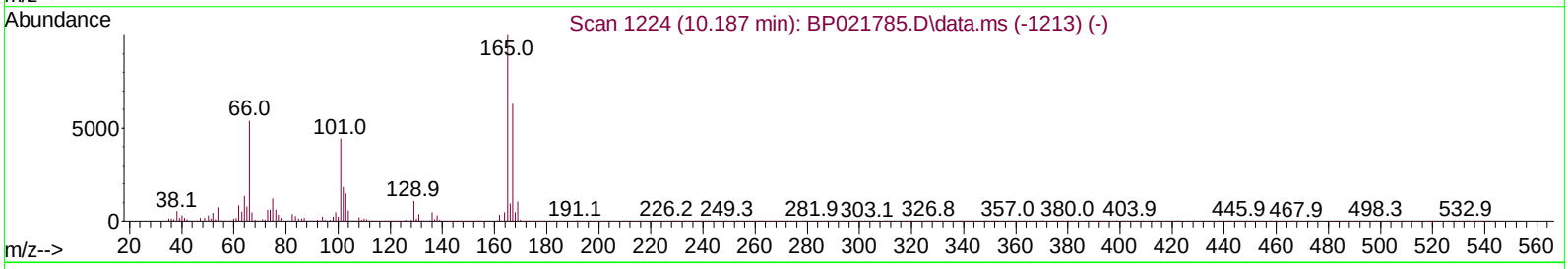
Instrument :
BNA_P
ClientSampleId :
C0SX1

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 09/04/2024
Supervised By :mohammad ahmed 09/07/2024







TIC: BP021798.D\data.ms

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

10.187min (-0.006) 22.98 ng/ul

response 418254

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.10	62.73
101.00	41.40	42.61
0.00	0.00	0.00

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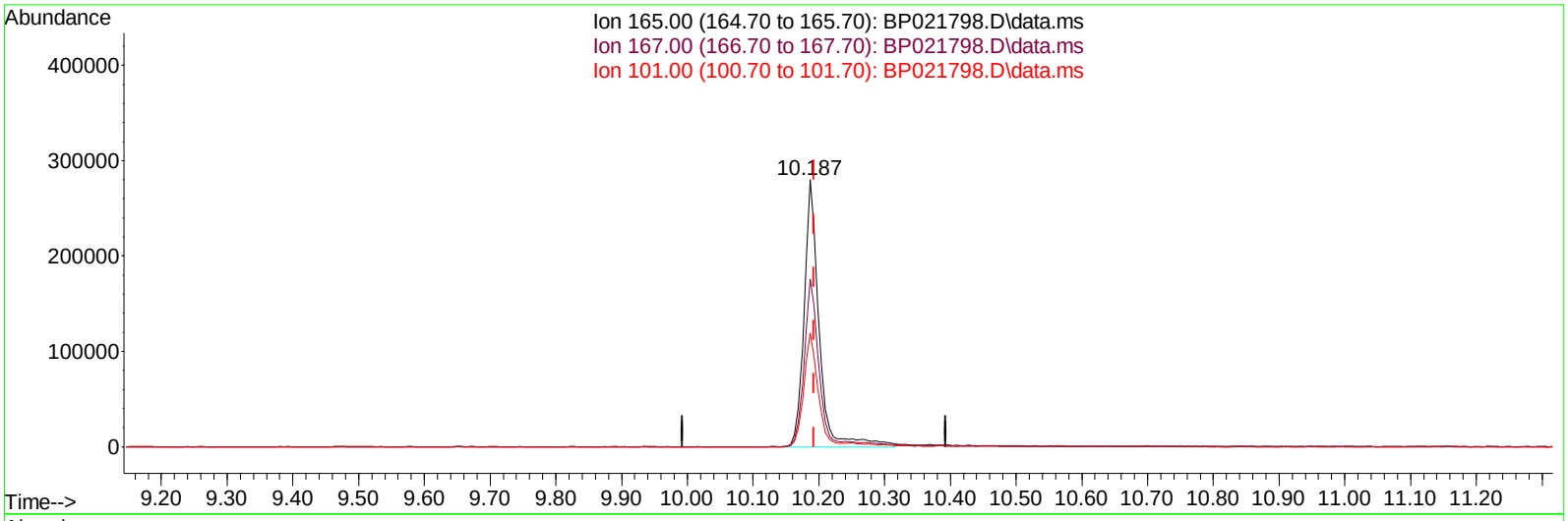

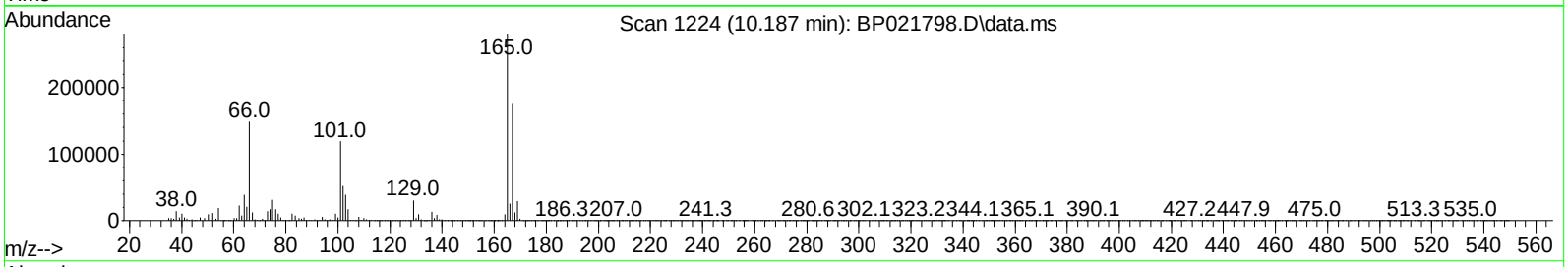

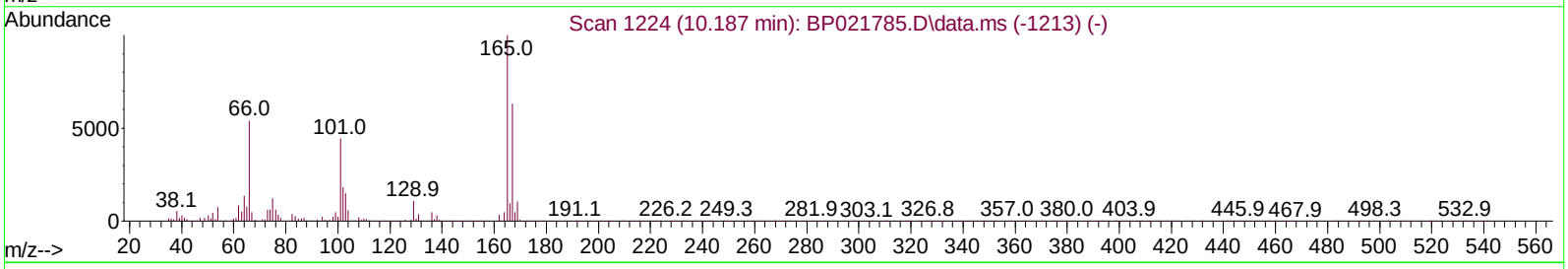

TIC: BP021798.D\data.ms

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

10.187min (-0.006) 24.67 ng/ul m

response 449030

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	64.10	62.73
101.00	41.40	42.61
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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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

Instrument :
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	284654	20.000	ng/ul	0.00
20) Naphthalene-d8	10.563	136	1130963	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.422	164	720983	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.233	188	1602317	20.000	ng/ul	0.00
79) Chrysene-d12	21.674	240	1587838	20.000	ng/ul	0.00
88) Perylene-d12	25.074	264	1760572	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	36971	4.312	ng/uL	0.00
4) Pyridine-d5	3.693	84	95352	3.845	ng/ul	0.00
7) Phenol-d5	6.952	99	268440	8.837	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.110	67	441477	26.704	ng/ul	0.00
11) 2-Chlorophenol-d4	7.322	132	485549	24.757	ng/ul	0.00
15) 4-Methylphenol-d8	8.481	113	407929	17.378	ng/ul	0.00
21) Nitrobenzene-d5	8.928	128	256089	28.145	ng/ul	0.00
24) 2-Nitrophenol-d4	9.651	143	253757	27.475	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.187	165	449030m	24.671	ng/ul	0.00
31) 4-Chloroaniline-d4	10.698	131	229899	8.682	ng/ul	0.00
46) Dimethylphthalate-d6	13.822	166	1546155	28.026	ng/ul	-0.02
49) Acenaphthylene-d8	14.110	160	1697249	27.402	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.634	143	64327	6.129	ng/ul	0.00
60) Fluorene-d10	15.428	176	1321866	28.501	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.557	200	207264	23.518	ng/ul	0.00
73) Anthracene-d10	17.339	188	2330649	30.649	ng/ul	0.00
81) Pyrene-d10	19.704	212	2788517	30.722	ng/ul	0.01
92) Benzo(a)pyrene-d12	24.845	264	2914240	31.042	ng/ul	0.00

Target Compounds Qvalue

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

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