Quantitation Report (Qedit)

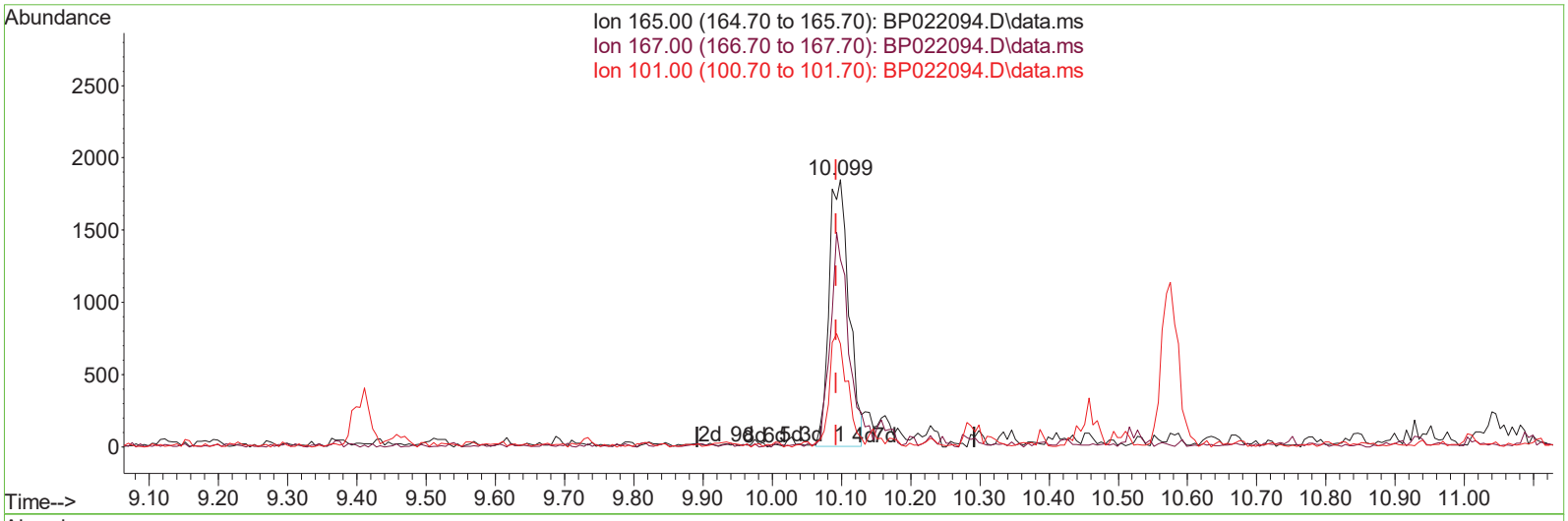
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022094.D
Acq On : 28 Sep 2024 06:32
Operator : RC/JU
Sample : P3910-15 10X
Misc :
ALS Vial : 23 Sample Multiplier: 1

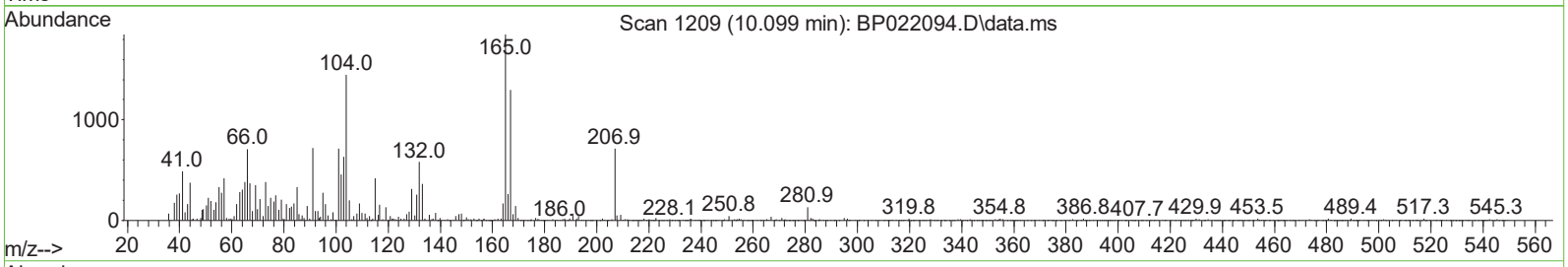
Quant Time: Sep 28 06:58:29 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 24 15:23:13 2024
Response via : Initial Calibration

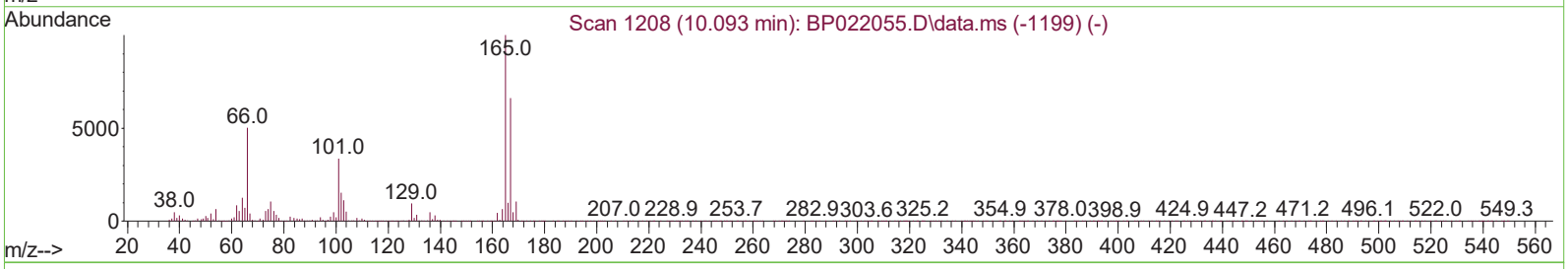
Instrument :
BNA_P
ClientSampleId :
DDC35

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 10/01/2024
Supervised By :mohammad ahmed 10/01/2024







TIC: BP022094.D\data.ms

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

10.099min (+ 0.006) 0.68 ng/ul

response 3677

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	62.40	70.24
101.00	33.20	38.64
0.00	0.00	0.00

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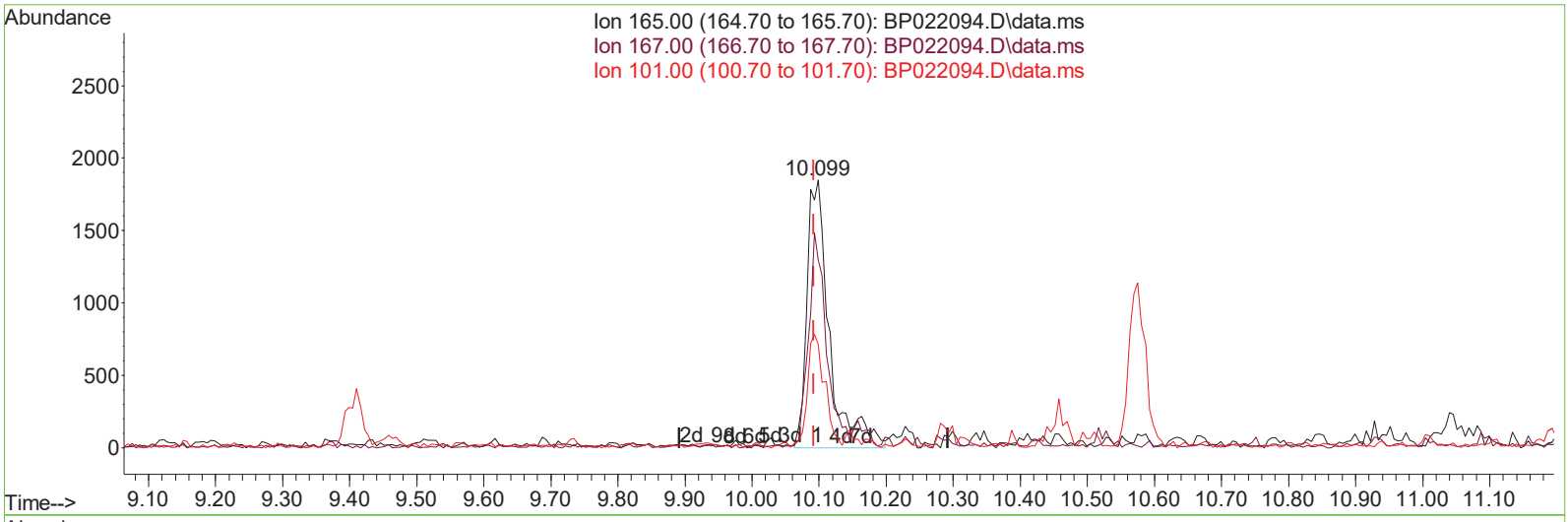

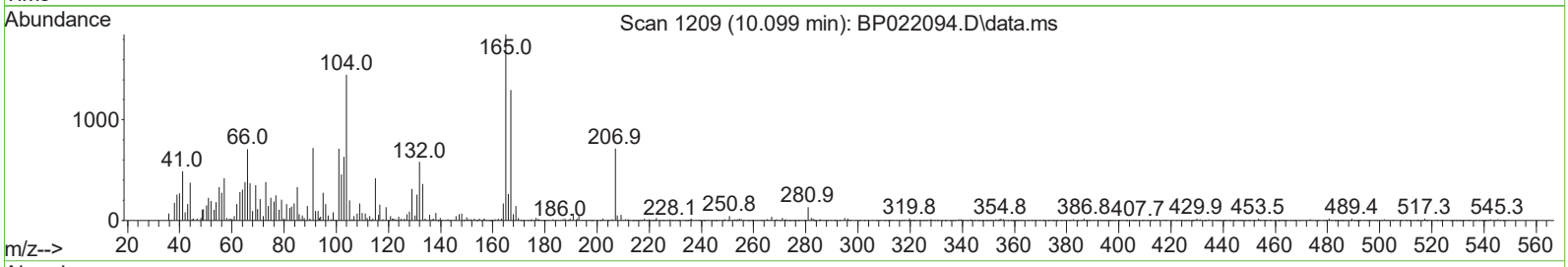

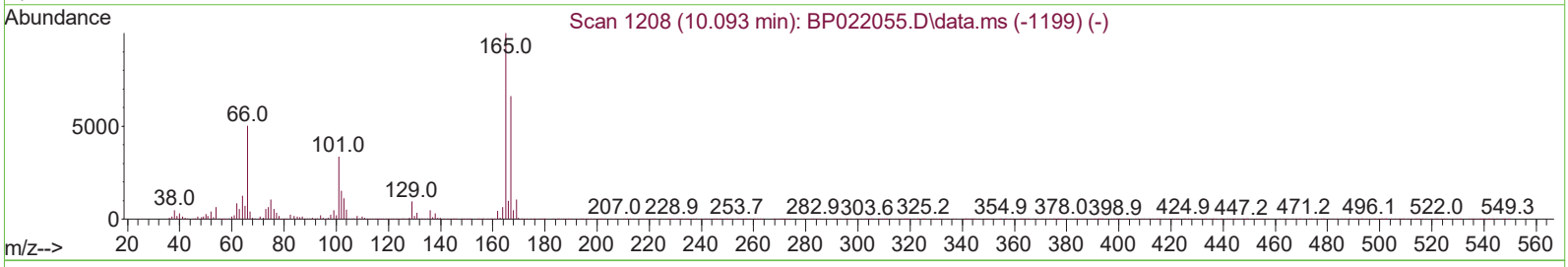

TIC: BP022094.D\data.ms

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

10.099min (+ 0.006) 0.80 ng/ul m

response 4295

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	62.40	70.24
101.00	33.20	38.64
0.00	0.00	0.00

Quantitation Report (LSC Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092724\
Data File : BP022094.D
Acq On : 28 Sep 2024 06:32
Operator : RC/JU
Sample : P3910-15 10X
Misc :
ALS Vial : 23 Sample Multiplier: 1

Quant Time: Sep 28 06:59:41 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 24 15:23:13 2024
Response via : Initial Calibration

Instrument :
BNA_P
ClientSampleId :
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Manual IntegrationsAPPROVED

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Supervised By :mohammad ahmed 10/01/2024

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.705	152	77800	20.000	ng/ul	0.00
20) Naphthalene-d8	10.457	136	293027	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.322	164	225333	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.122	188	533888	20.000	ng/ul	0.00
79) Chrysene-d12	21.539	240	668329	20.000	ng/ul	0.00
88) Perylene-d12	24.827	264	795447	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.223	96	908	0.457	ng/uL	0.00
4) Pyridine-d5	3.646	84	11129	1.960	ng/ul	0.01
7) Phenol-d5	6.887	99	9696	1.524	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.040	67	9487	2.421	ng/ul	0.00
11) 2-Chlorophenol-d4	7.246	132	4525	0.983	ng/ul	0.00
15) 4-Methylphenol-d8	8.405	113	10986	2.178	ng/ul	0.00
21) Nitrobenzene-d5	8.840	128	5618	2.562	ng/ul	0.00
24) 2-Nitrophenol-d4	9.563	143	1435	0.577	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.099	165	4295m	0.795	ng/ul	0.00
31) 4-Chloroaniline-d4	10.599	131	18107	2.810	ng/ul	0.00
46) Dimethylphthalate-d6	13.722	166	58395	3.352	ng/ul	0.00
49) Acenaphthylene-d8	14.004	160	57311	3.108	ng/ul	0.00
54) 4-Nitrophenol-d4	14.528	143	955	0.321	ng/ul	0.00
60) Fluorene-d10	15.328	176	54511	3.538	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.451	200	839	0.253	ng/ul	0.00
73) Anthracene-d10	17.210	188	95162	3.818	ng/ul	0.00
81) Pyrene-d10	19.592	212	125870	3.635	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.598	264	150266	3.592	ng/ul	-0.01
Target Compounds						Qvalue
36) 2-Methylnaphthalene	12.116	142	13051	1.139	ng/ul	95

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

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