Quantitation Report (Qedit)

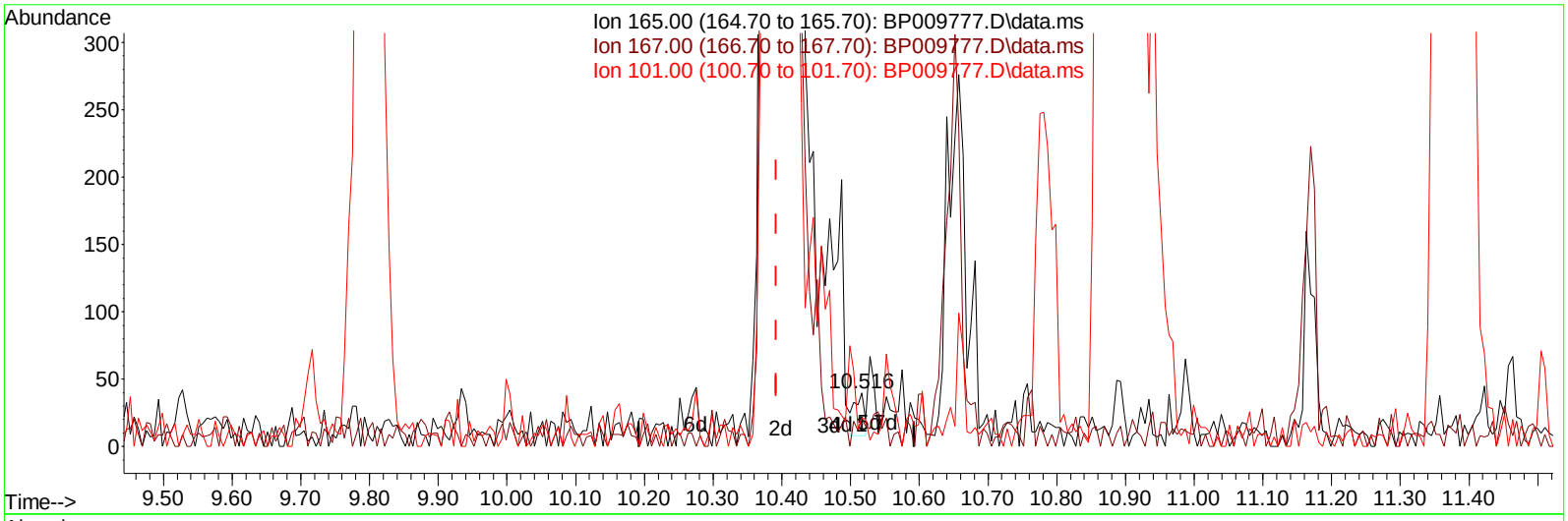
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
Data File : BP009777.D
Acq On : 12 Apr 2022 10:34
Operator : CG/JU
Sample : N2338-02DL 10X
Misc :
ALS Vial : 35 Sample Multiplier: 1

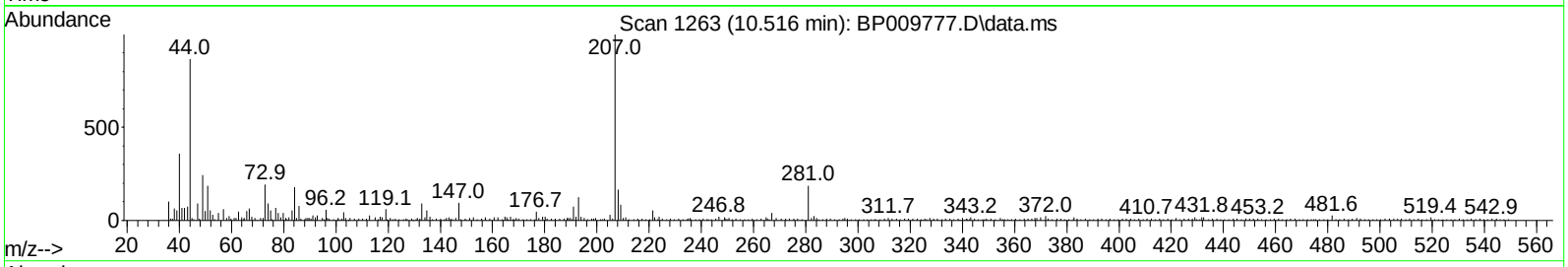
Quant Time: Apr 13 02:19:33 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Apr 13 02:15:22 2022
Response via : Initial Calibration

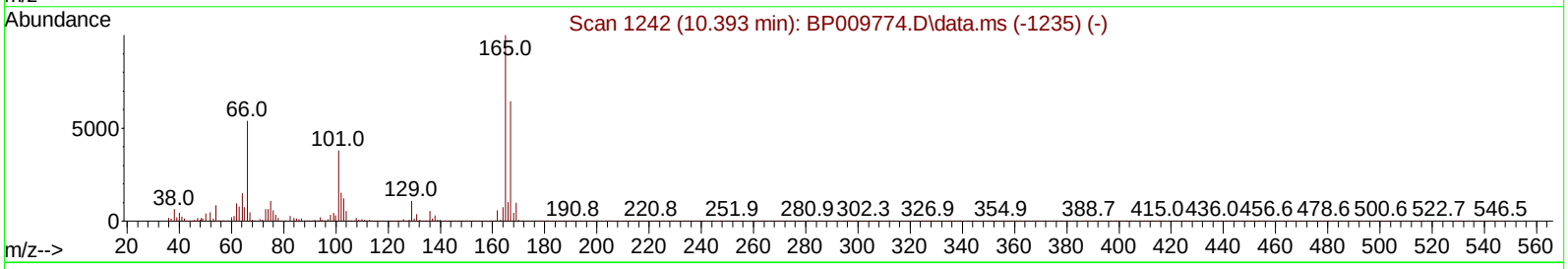
Instrument :
BNA_P
ClientSampleId :
C0J71DL

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/13/2022
Supervised By :Yogesh Patel 04/13/2022







TIC: BP009777.D\data.ms

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

10.516min (+ 0.124) 0.01 ng/ul

response 32

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	52.50
101.00	49.80	52.50
0.00	0.00	0.00

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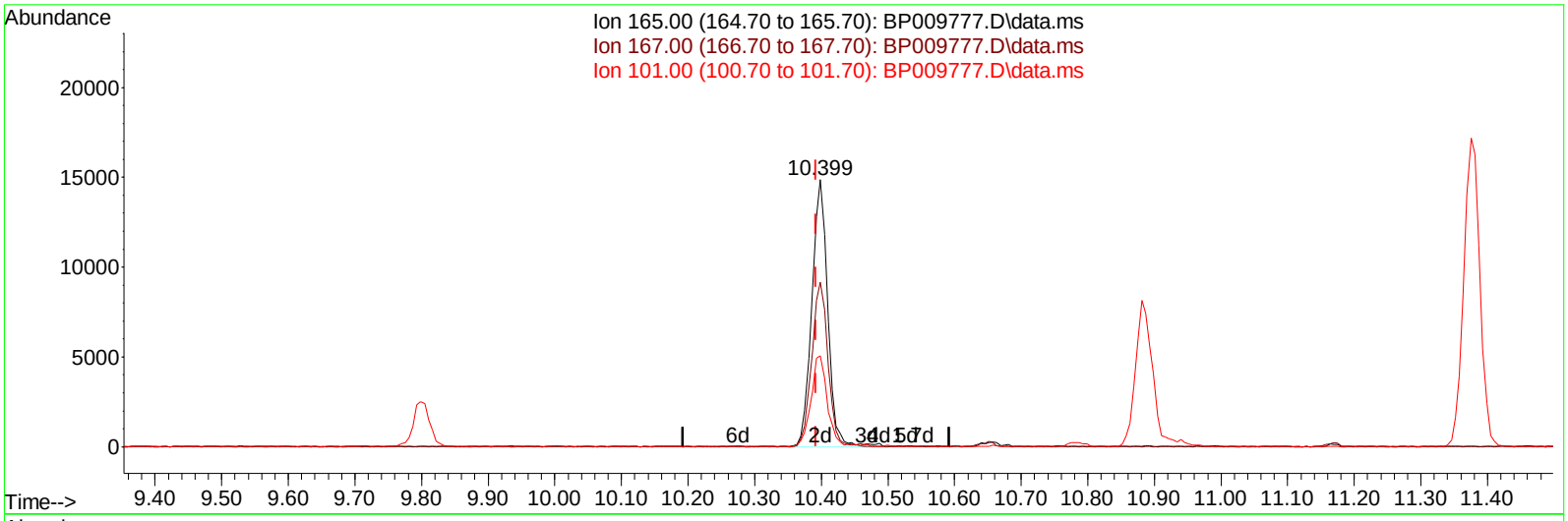

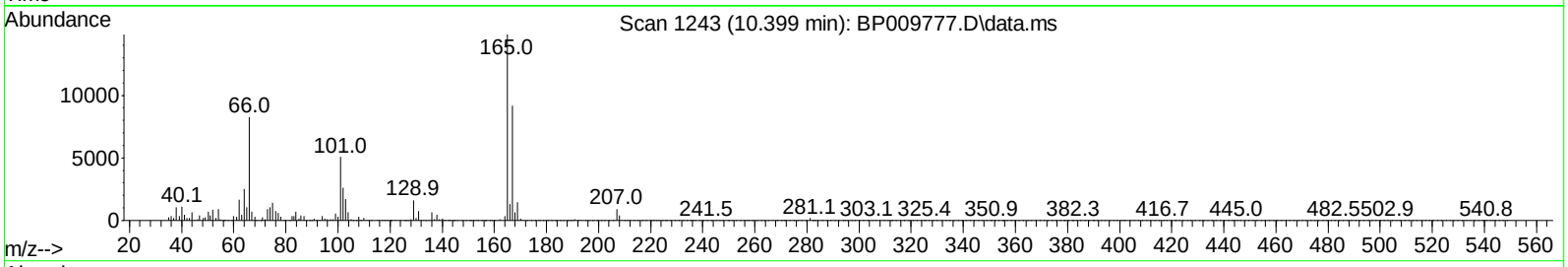

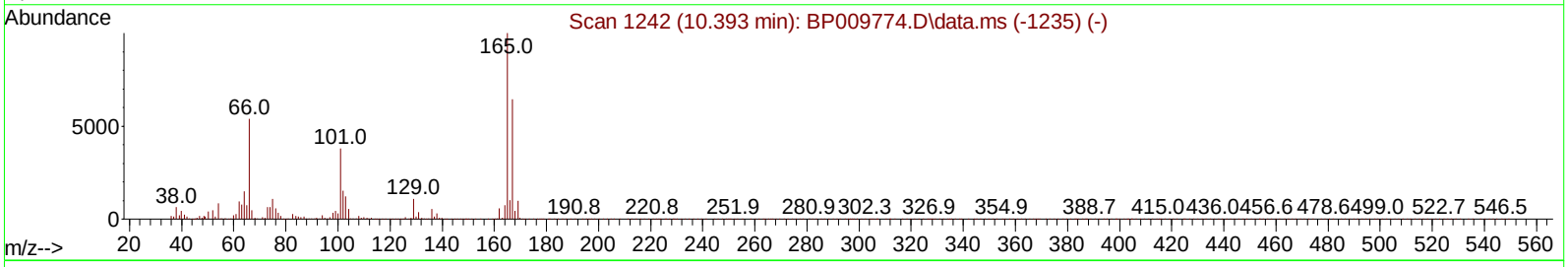

TIC: BP009777.D\data.ms

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

10.399min (+ 0.006) 3.94 ng/ul m

response 24646

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	61.60
101.00	49.80	34.20#
0.00	0.00	0.00

Quantitation Report (QT/LSC Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041122\
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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/13/2022
Supervised By :Yogesh Patel 04/13/2022

Compound	R.T.	QIon	Response	Conc	Units		Dev(Min)
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.981	152	88426	20.000	ng/ul		0.01
20) Naphthalene-d8	10.787	136	378071	20.000	ng/ul		0.01
38) Acenaphthene-d10	14.604	164	272917	20.000	ng/ul		0.00
64) Phenanthrene-d10	17.351	188	621319	20.000	ng/ul	#	0.00
79) Chrysene-d12	21.433	240	664112	20.000	ng/ul		0.00
88) Perylene-d12	23.904	264	675956	20.000	ng/ul		0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	0.000	96	0d	0.000	ng/uL		
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul		
7) Phenol-d5	7.122	99	7616	0.981	ng/ul		0.00
9) Bis-(2-Chloroethyl)eth...	7.293	67	42838	9.269	ng/ul		0.00
11) 2-Chlorophenol-d4	7.505	132	21466	3.666	ng/ul		0.00
15) 4-Methylphenol-d8	8.675	113	13903	2.145	ng/ul		0.00
21) Nitrobenzene-d5	9.134	128	13231	4.853	ng/ul		0.00
24) 2-Nitrophenol-d4	9.858	143	13859	4.637	ng/ul		0.00
28) 2,4-Dichlorophenol-d3	10.399	165	24646m	3.940	ng/ul		0.00
31) 4-Chloroaniline-d4	10.910	131	36731	4.123	ng/ul		0.00
46) Dimethylphthalate-d6	14.010	166	108167	5.046	ng/ul		0.00
49) Acenaphthylene-d8	14.298	160	112174	4.417	ng/ul		0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul		
60) Fluorene-d10	15.593	176	94649	5.248	ng/ul		0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	10233	2.780	ng/ul		0.00
73) Anthracene-d10	17.451	188	151951	5.108	ng/ul		0.00
81) Pyrene-d10	19.681	212	189936	5.237	ng/ul		0.00
92) Benzo(a)pyrene-d12	23.739	264	172988	4.914	ng/ul		0.00
Target Compounds							Qvalue
8) Phenol	7.152	94	14856	1.912	ng/ul	#	82
22) Nitrobenzene	9.181	77	2137323	305.418	ng/ul	#	84
32) 4-Chloroaniline	10.934	127	15459	1.754	ng/ul		94
39) 1,2,4,5-Tetrachloroben...	12.804	216	40480	4.848	ng/ul		98
75) 1,2,3,4-Tetrachloroben...	13.404	216	70791	7.881	ng/uL	#	86

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

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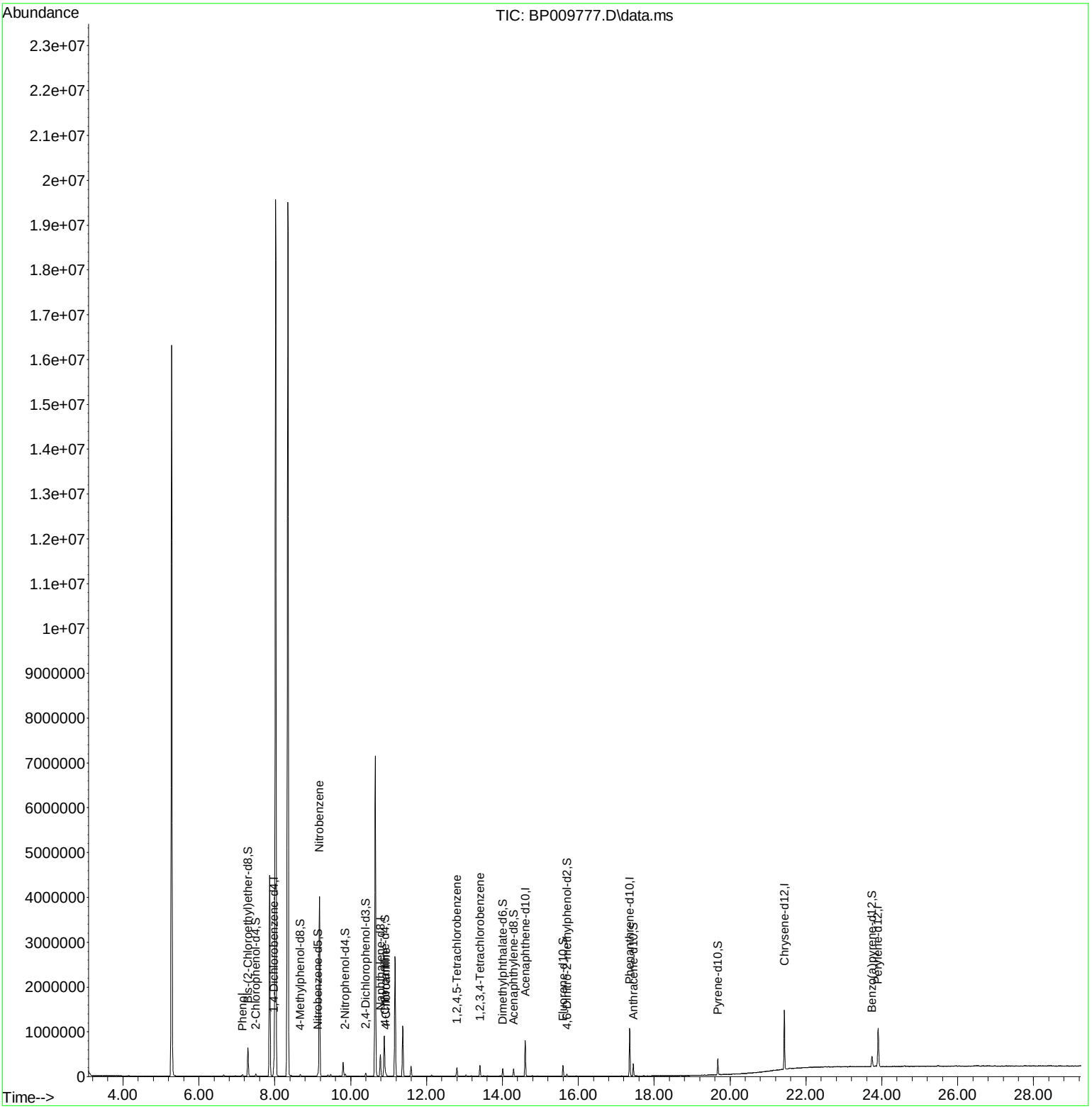
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