

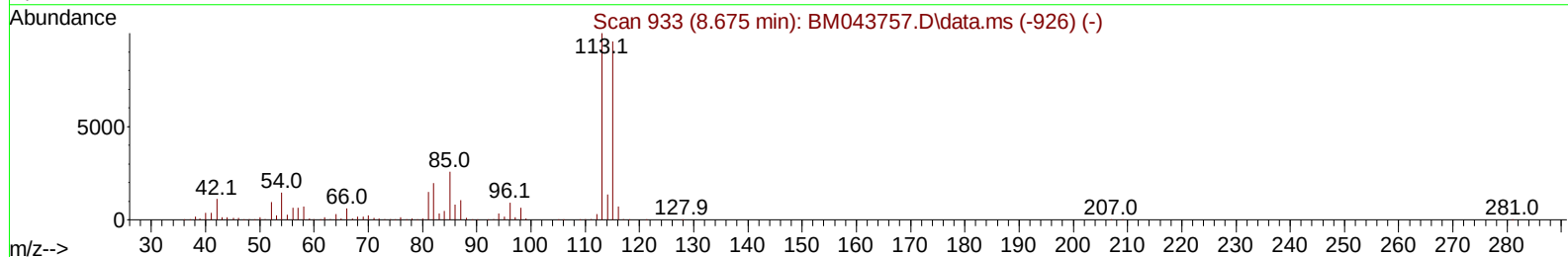
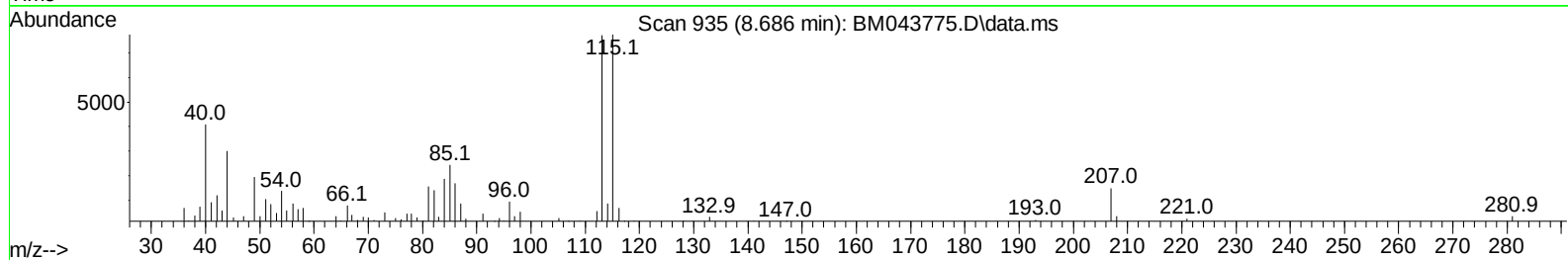
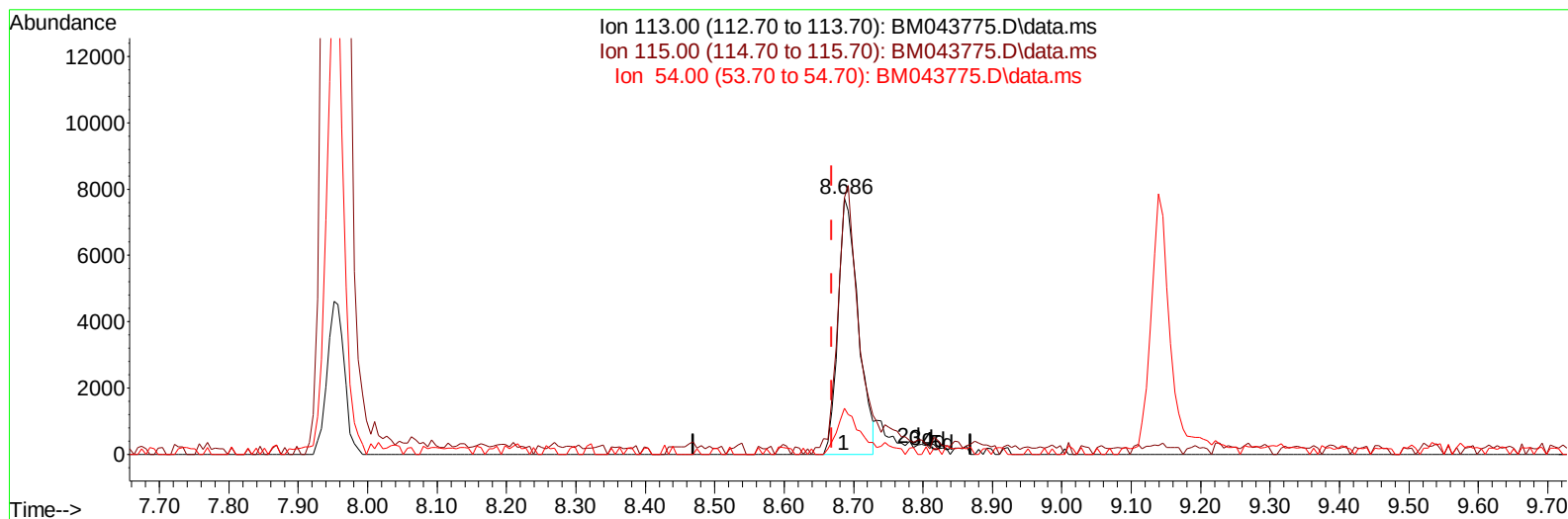
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
Data File : BM043775.D
Acq On : 06 Jan 2024 00:42
Operator : MA/JU
Sample : 05919-14MEDL 20X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCF46MEDL

Manual IntegrationsAPPROVED

Quant Time: Jan 06 01:22:29 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Dec 29 22:31:57 2023
Response via : Initial Calibration

Reviewed By :Yogesh Patel 01/08/2024
Supervised By :mohammad ahmed 01/08/2024



TIC: BM043775.D\data.ms

(15) 4-Methylphenol-d8 (S)

8.686min (+ 0.017) 0.63 ng/u1

response 15589

Ion	Exp%	Act%
113.00	100.00	100.00
115.00	95.60	100.27
54.00	15.00	17.96
0.00	0.00	0.00

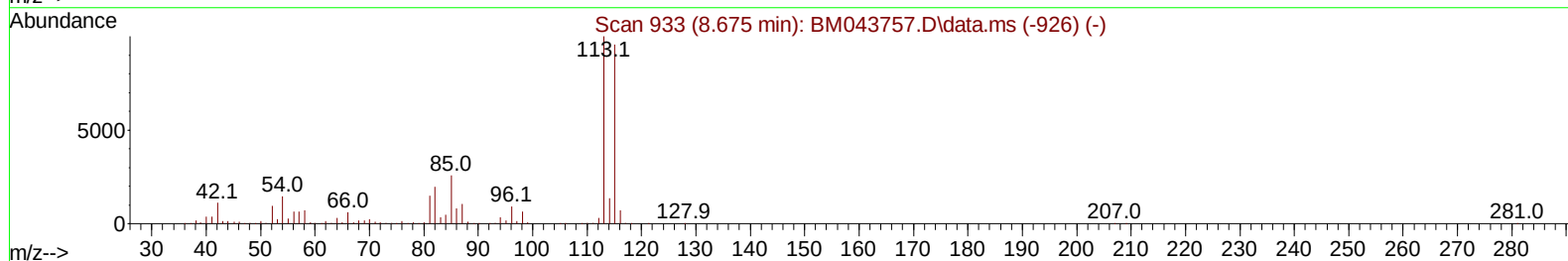
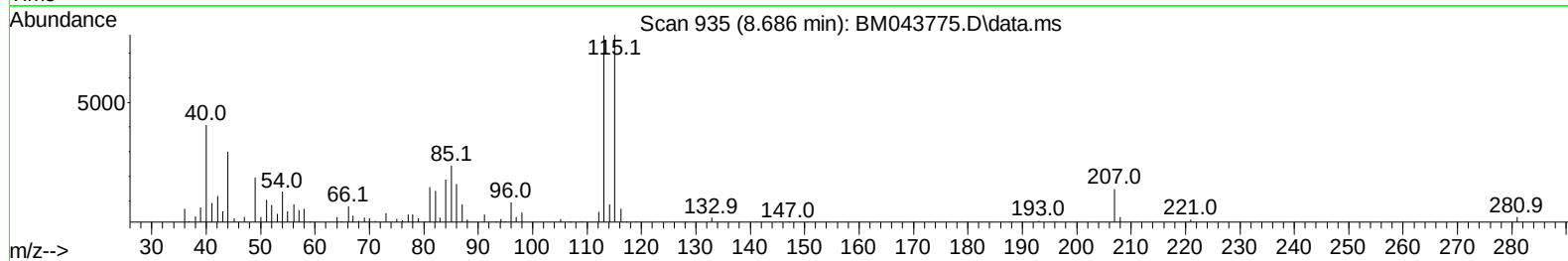
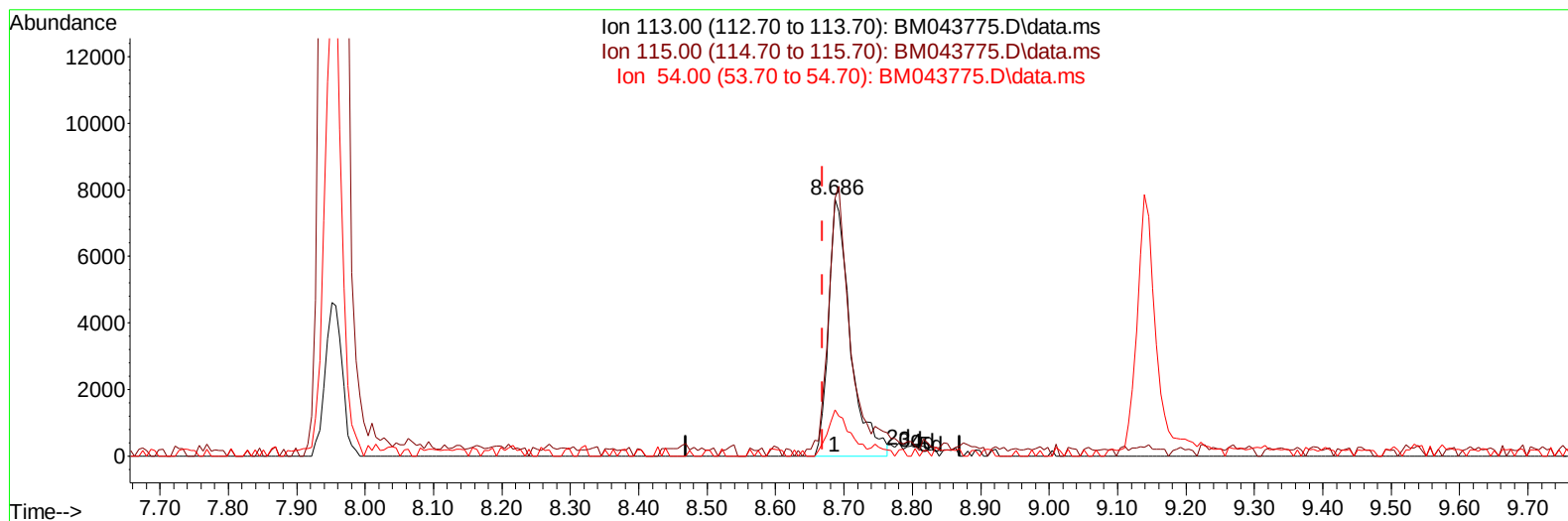
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
Data File : BM043775.D
Acq On : 06 Jan 2024 00:42
Operator : MA/JU
Sample : 05919-14MEDL 20X
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GCF46MEDL

Manual IntegrationsAPPROVED

Quant Time: Jan 06 02:53:31 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Dec 29 22:31:57 2023
Response via : Initial Calibration

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Supervised By :mohammad ahmed 01/08/2024



TIC: BM043775.D\data.ms

(15) 4-Methylphenol-d8 (S)

8.686min (+ 0.017) 0.69 ng/ul m

response 16997

Ion	Exp%	Act%
113.00	100.00	100.00
115.00	95.60	100.27
54.00	15.00	17.96
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
 Data File : BM043775.D
 Acq On : 06 Jan 2024 00:42
 Operator : MA/JU
 Sample : 05919-14MEDL 20X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :

BNA_M

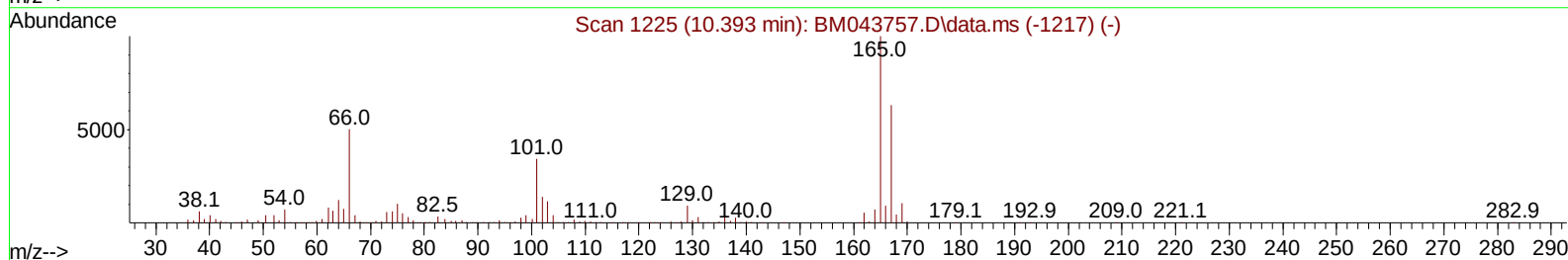
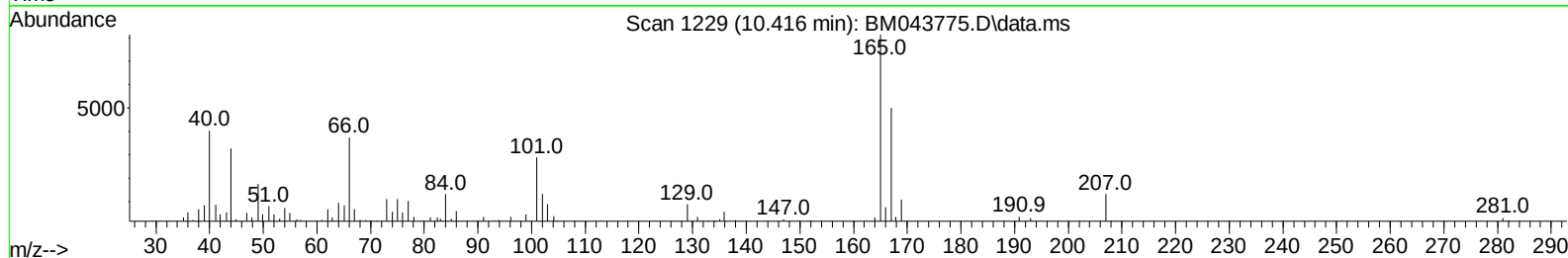
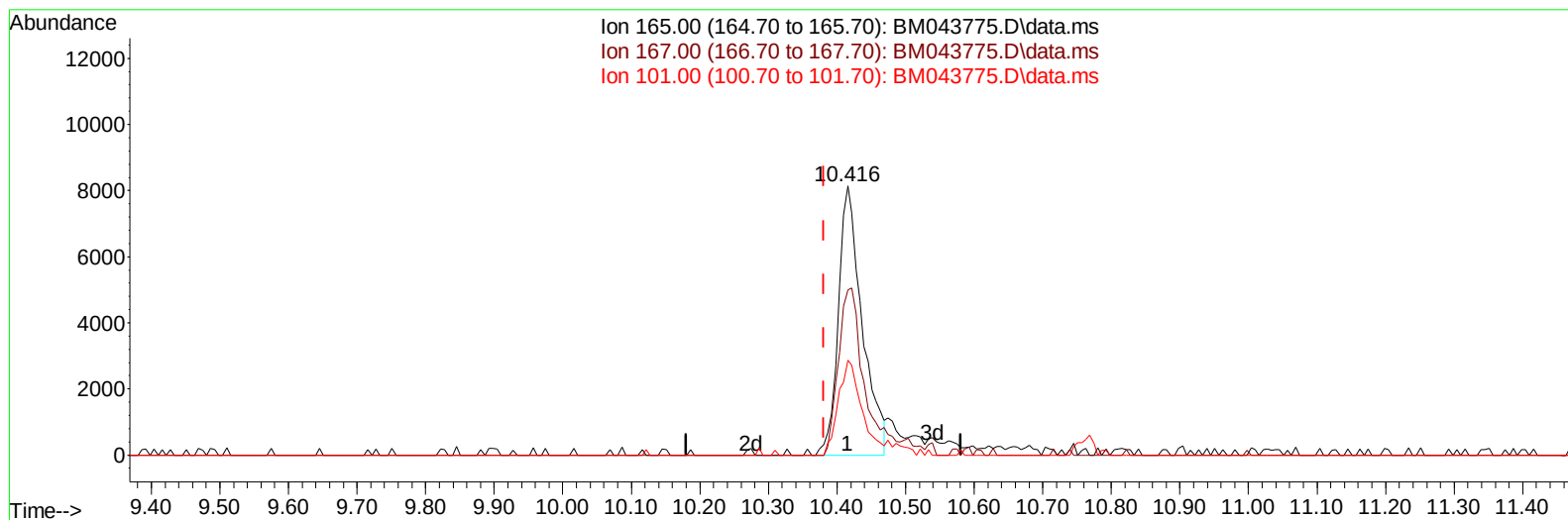
ClientSampleId :

GCF46MEDL

Manual IntegrationsAPPROVED

Quant Time: Jan 06 01:23:00 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 29 22:31:57 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 01/08/2024
 Supervised By :mohammad ahmed 01/08/2024



TIC: BM043775.D\data.ms

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

10.416min (+ 0.035) 1.01 ng/ul

response 19545

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	61.80	61.40
101.00	37.00	35.33
0.00	0.00	0.00

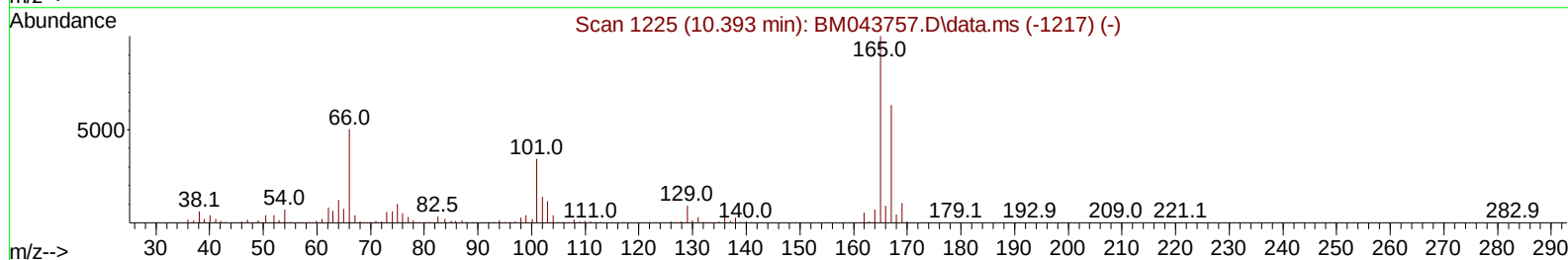
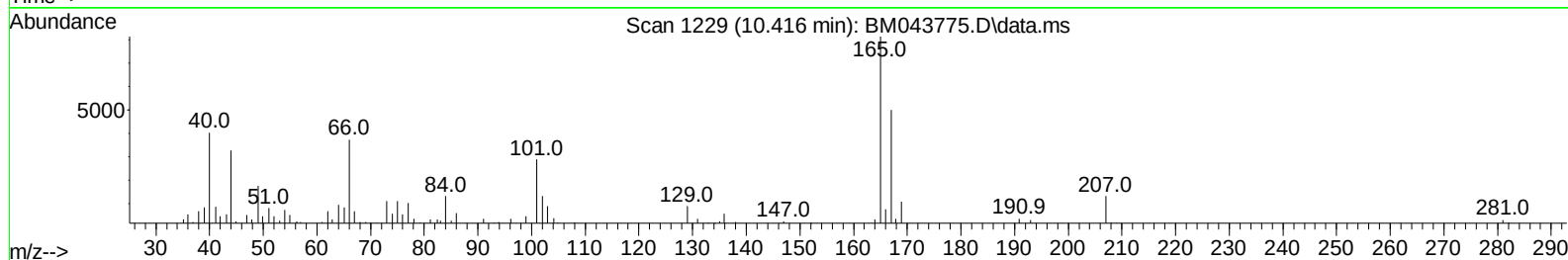
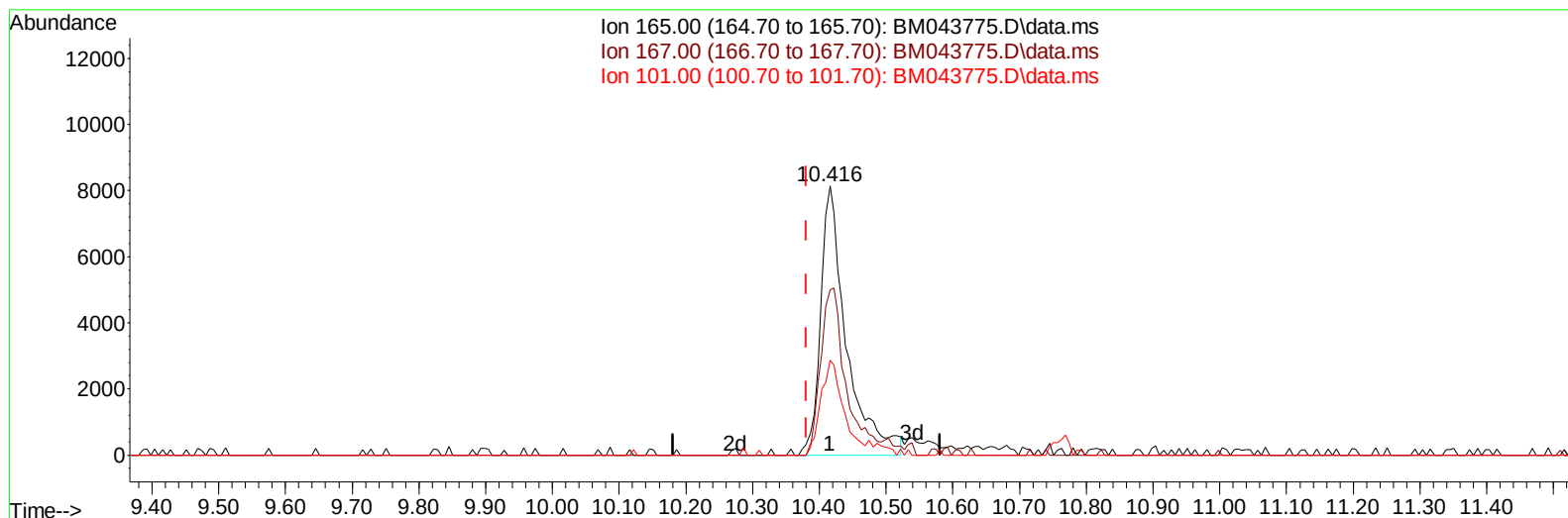
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010524\
 Data File : BM043775.D
 Acq On : 06 Jan 2024 00:42
 Operator : MA/JU
 Sample : 05919-14MEDL 20X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 GCF46MEDL

Manual IntegrationsAPPROVED

Quant Time: Jan 06 01:23:00 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 29 22:31:57 2023
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Reviewed By :Yogesh Patel 01/08/2024
 Supervised By :mohammad ahmed 01/08/2024



TIC: BM043775.D\data.ms

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

10.416min (+ 0.035) 1.12 ng/ul m

response 21752

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	61.80	61.40
101.00	37.00	35.33
0.00	0.00	0.00

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 Sample : 05919-14MEDL 20X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
 GCF46MEDL

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 01/08/2024
 Supervised By :mohammad ahmed 01/08/2024

Quant Time: Jan 06 02:53:31 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.957	152	279796	20.000	ng/ul	0.00
20) Naphthalene-d8	10.763	136	1098232	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.598	164	634225	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.351	188	1292405	20.000	ng/ul	0.01
79) Chrysene-d12	21.533	240	1166687	20.000	ng/ul	0.01
88) Perylene-d12	23.950	264	1365292	20.000	ng/ul	0.02
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	7.139	99	29826	0.937	ng/ul	0.02
9) Bis-(2-Chloroethyl)eth...	7.298	67	20534	1.019	ng/ul	0.00
11) 2-Chlorophenol-d4	7.486	132	24716	1.017	ng/ul	0.00
15) 4-Methylphenol-d8	8.686	113	16997m	0.688	ng/ul	0.02
21) Nitrobenzene-d5	9.139	128	11116	1.046	ng/ul	0.01
24) 2-Nitrophenol-d4	9.857	143	12412	1.131	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.416	165	21752m	1.122	ng/ul	0.03
31) 4-Chloroaniline-d4	0.000	131	0d	0.000	ng/ul	
46) Dimethylphthalate-d6	14.021	166	74747	1.309	ng/ul	0.00
49) Acenaphthylene-d8	14.292	160	85231	1.258	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.592	176	67002	1.419	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.739	200	6943	0.812	ng/ul	0.03
73) Anthracene-d10	17.445	188	100973	1.393	ng/ul	0.00
81) Pyrene-d10	19.739	212	112219	1.218	ng/ul	0.01
92) Benzo(a)pyrene-d12	23.791	264	106754	1.267	ng/ul	0.02
Target Compounds						
					Qvalue	
58) 2,3,4,6-Tetrachlorophenol	15.233	232	67007	5.997	ng/ul	92
71) Pentachlorophenol	17.003	266	2155841	203.288	ng/ul	99
72) Phenanthrene	17.392	178	146214	1.669	ng/ul	98

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

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