Quantitation Report (Qedit)

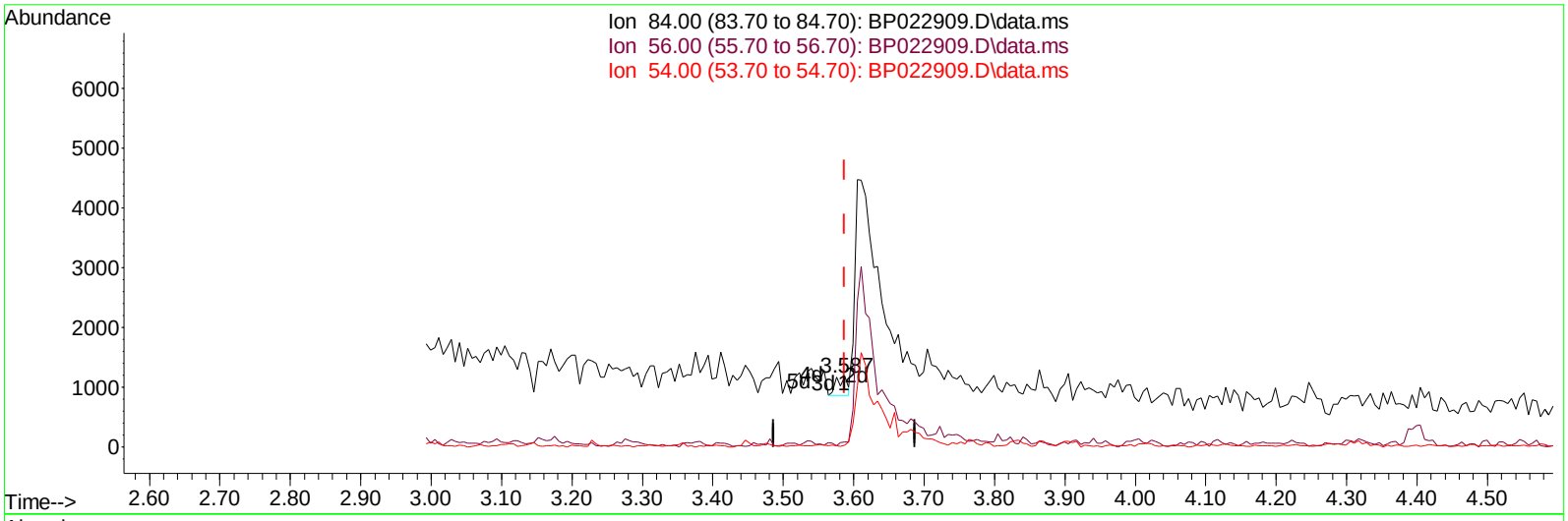
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022909.D
Acq On : 07 Nov 2024 01:24
Operator : RC/JU
Sample : P4622-10
Misc :
ALS Vial : 56 Sample Multiplier: 1

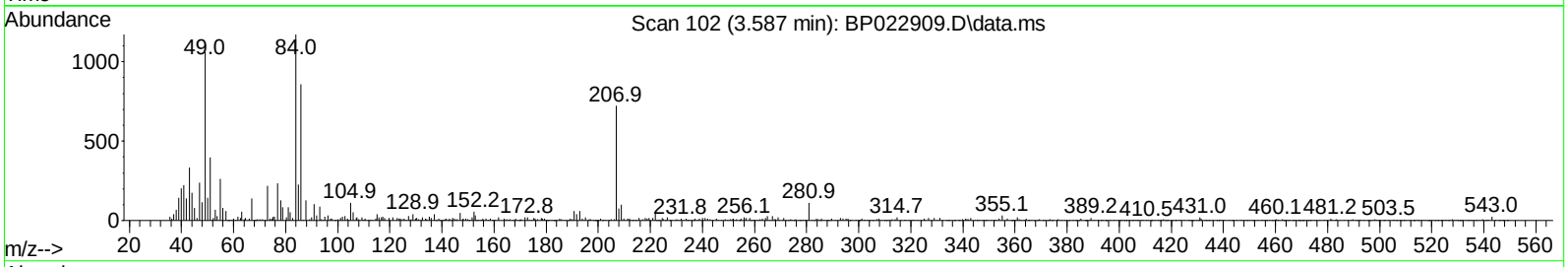
Quant Time: Nov 07 02:10:07 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 06 00:31:03 2024
Response via : Initial Calibration

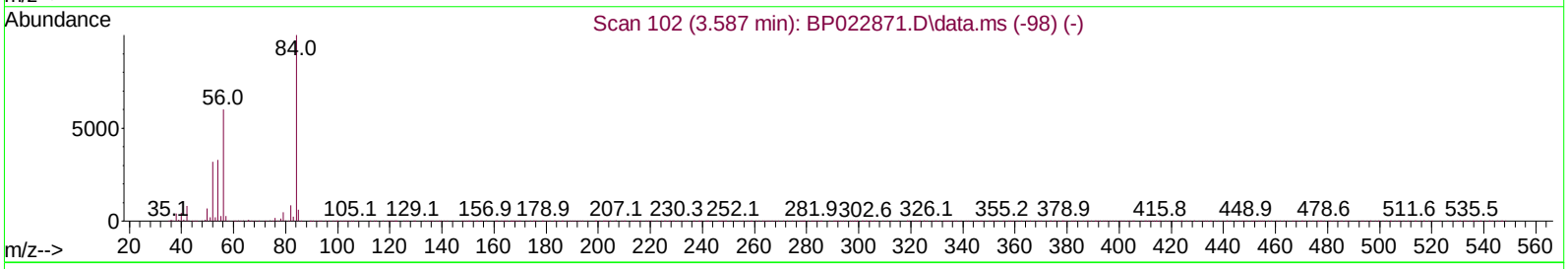
Instrument :
BNA_P
ClientSampleId :
BH6N6

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/07/2024
Supervised By :mohammad ahmed 11/08/2024







TIC: BP022909.D\data.ms

(4) Pyridine-d5 (S)

3.587min (+ 0.000) 0.05 ng/ul

response 334

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	61.10	6.91#
54.00	30.90	2.30#
0.00	0.00	0.00

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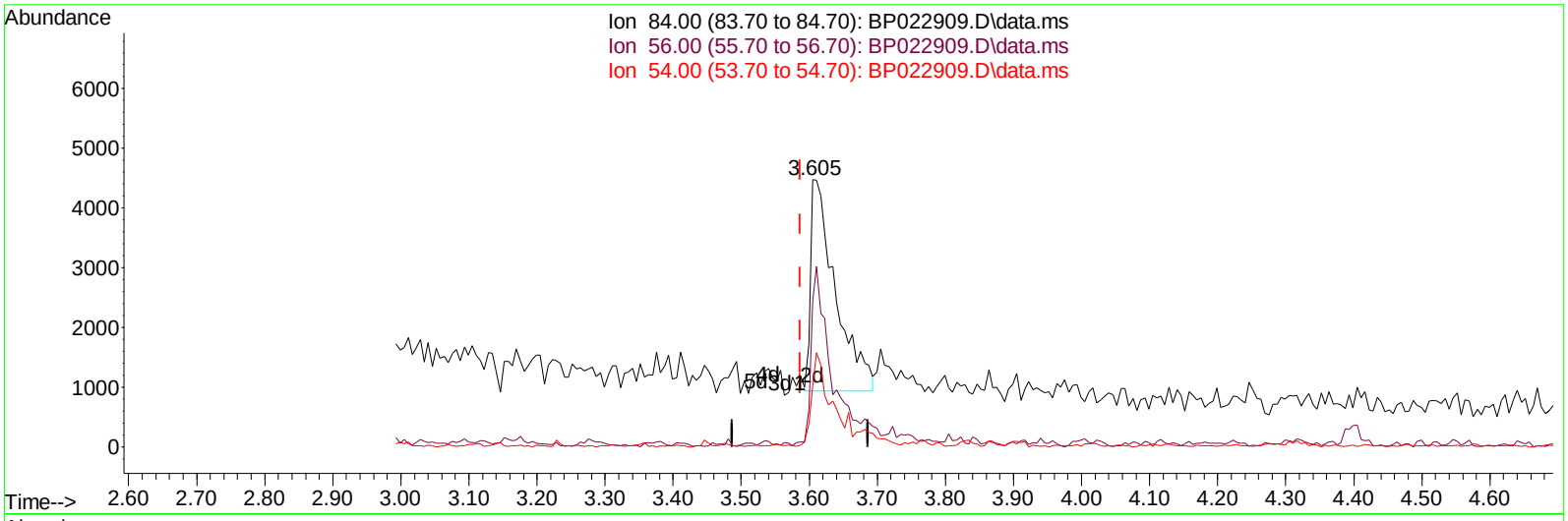

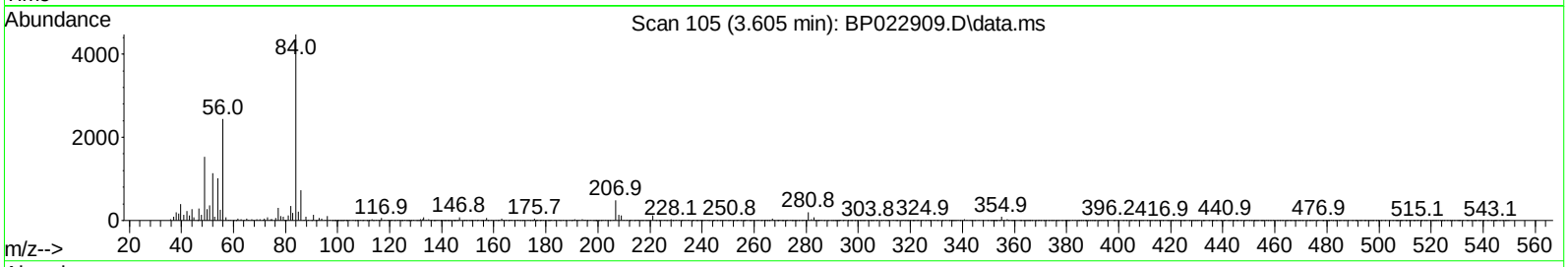

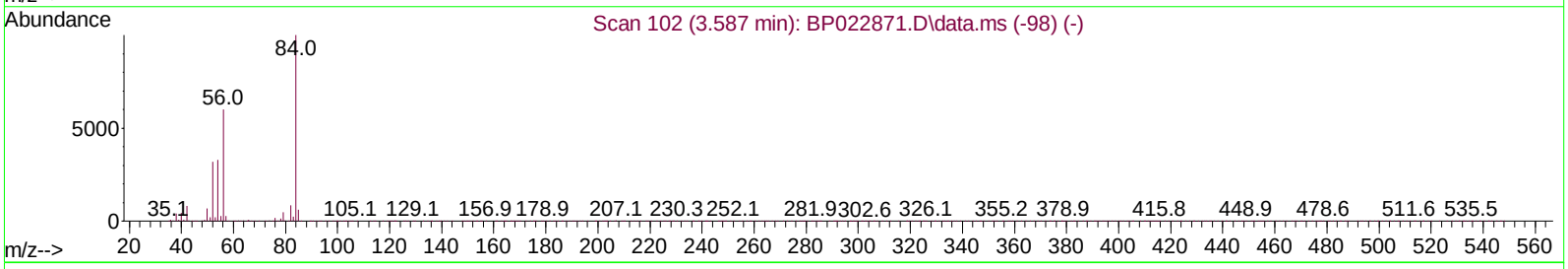

TIC: BP022909.D\data.ms

(4) Pyridine-d5 (S)

3.605min (+ 0.018) 1.30 ng/ul m

response 9054

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	61.10	54.49
54.00	30.90	22.85#
0.00	0.00	0.00

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QLast Update : Wed Nov 06 00:31:03 2024
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Instrument :
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ClientSampleId :
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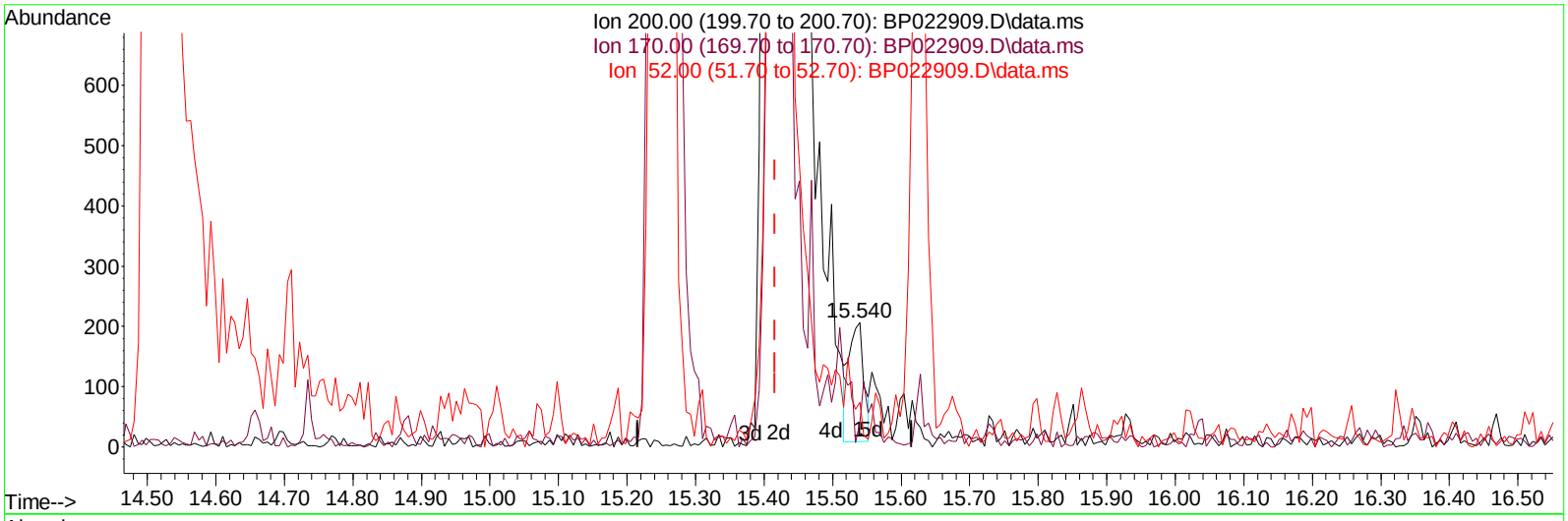

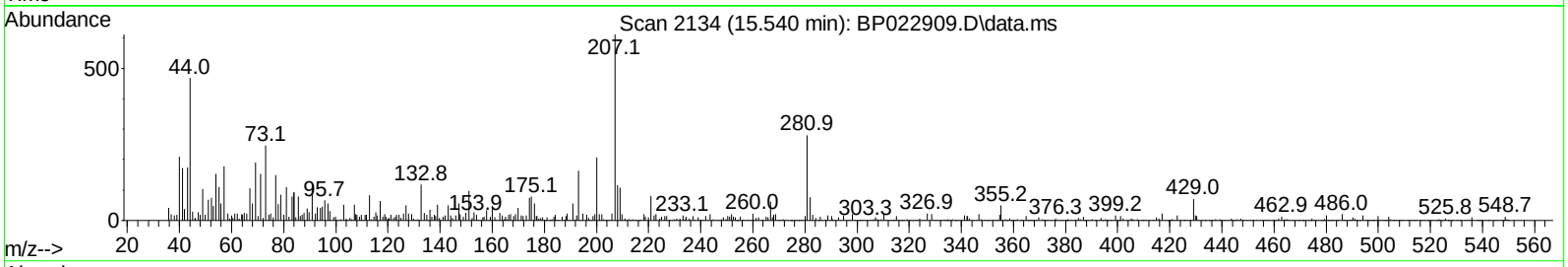

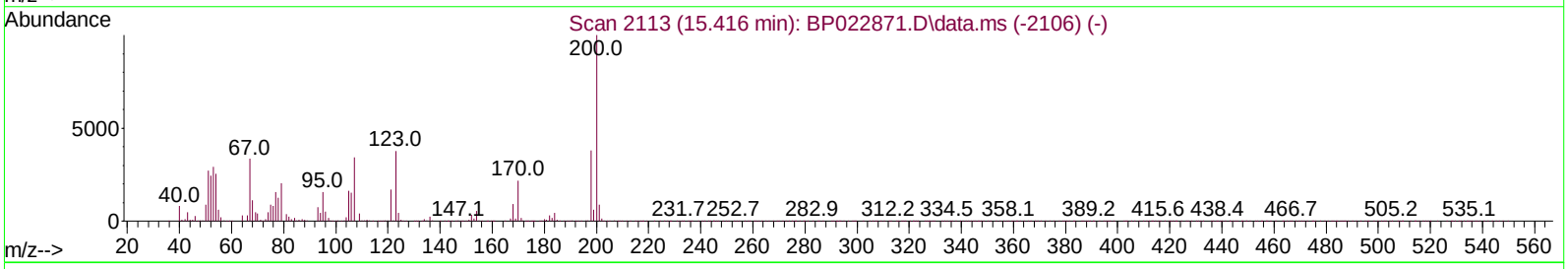

TIC: BP022909.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.540min (+ 0.124) 0.07 ng/ul

response 300

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.00	19.81
52.00	36.60	35.75
0.00	0.00	0.00

Quantitation Report (Qedit)

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ClientSampleId :
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Manual IntegrationsAPPROVED

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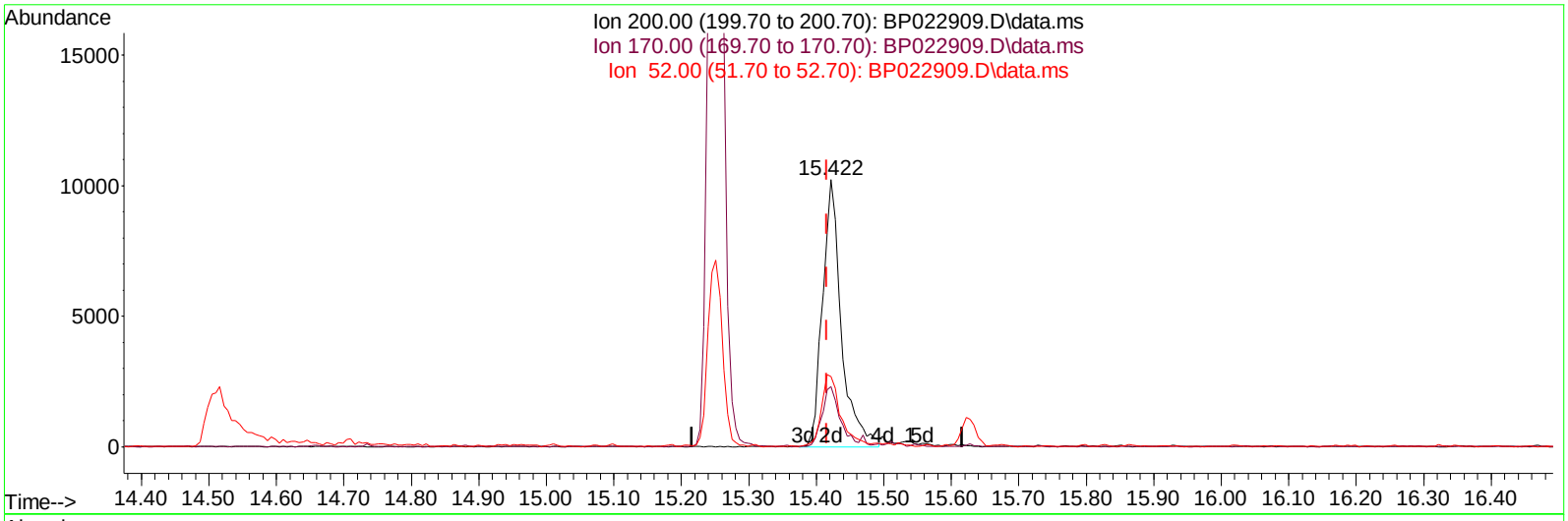

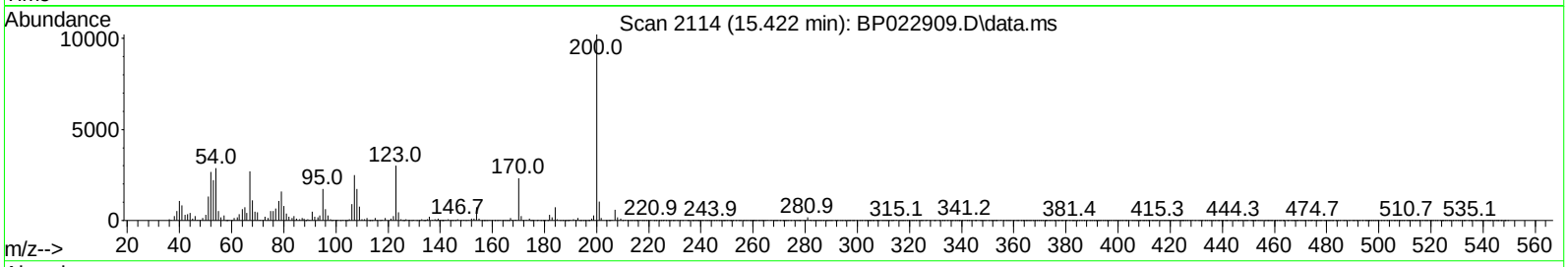

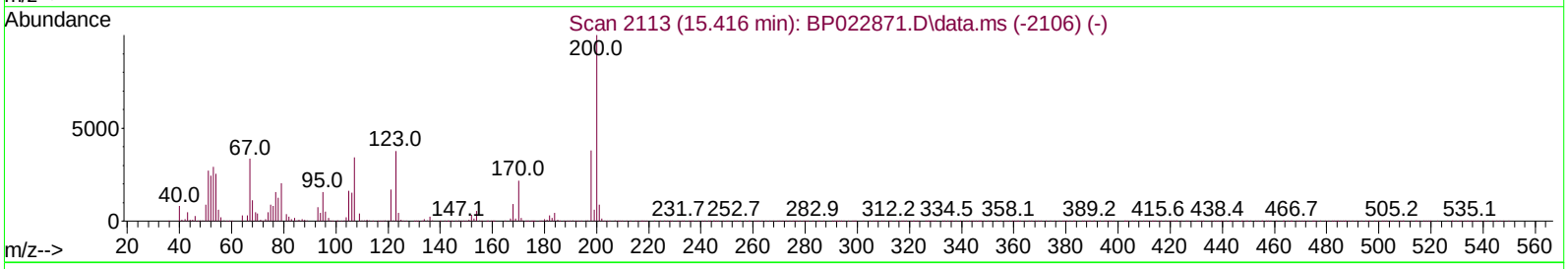

TIC: BP022909.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.422min (+ 0.006) 4.30 ng/ul m

response 19772

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.00	22.55
52.00	36.60	26.20#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/07/2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.622	152	98954	20.000	ng/ul	0.00
20) Naphthalene-d8	10.381	136	378859	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.240	164	292225	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.045	188	698901	20.000	ng/ul	0.00
79) Chrysene-d12	21.498	240	727556	20.000	ng/ul	# 0.00
88) Perylene-d12	24.763	264	866870	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	5850	2.297	ng/uL	0.00
4) Pyridine-d5	3.605	84	9054m	1.295	ng/ul	0.02
7) Phenol-d5	6.822	99	130089	15.618	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.969	67	73358	14.984	ng/ul	0.00
11) 2-Chlorophenol-d4	7.169	132	98638	16.689	ng/ul	0.00
15) 4-Methylphenol-d8	8.328	113	107735	16.207	ng/ul	0.00
21) Nitrobenzene-d5	8.763	128	49451	16.976	ng/ul	0.00
24) 2-Nitrophenol-d4	9.481	143	57851	17.153	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.022	165	125444	17.501	ng/ul	0.00
31) 4-Chloroaniline-d4	10.522	131	135857	16.040	ng/ul	0.00
46) Dimethylphthalate-d6	13.651	166	396411	17.073	ng/ul	-0.01
49) Acenaphthylene-d8	13.928	160	480307	19.477	ng/ul	0.00
54) 4-Nitrophenol-d4	14.504	143	45508	11.683	ng/ul	0.00
60) Fluorene-d10	15.251	176	405494	20.135	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.422	200	19772m	4.301	ng/ul	0.00
73) Anthracene-d10	17.163	188	658706	20.035	ng/ul	0.00
81) Pyrene-d10	19.539	212	846798	22.009	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.533	264	950688	20.535	ng/ul	0.00
Target Compounds						Qvalue

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

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