

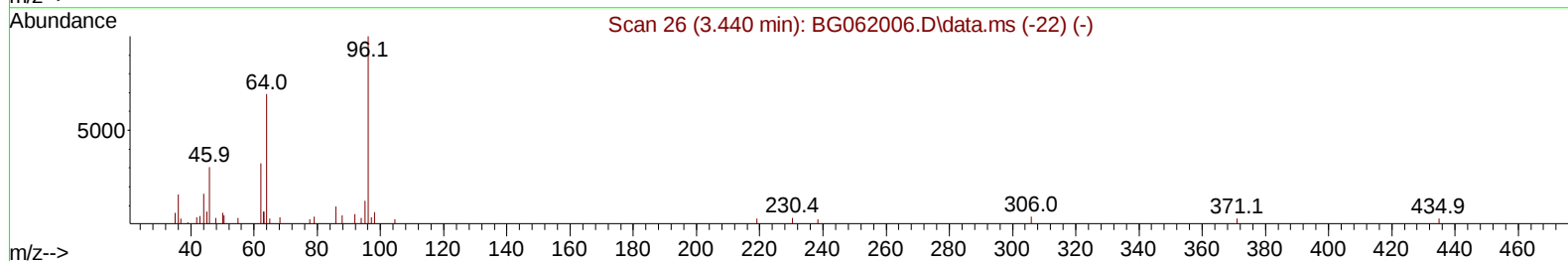
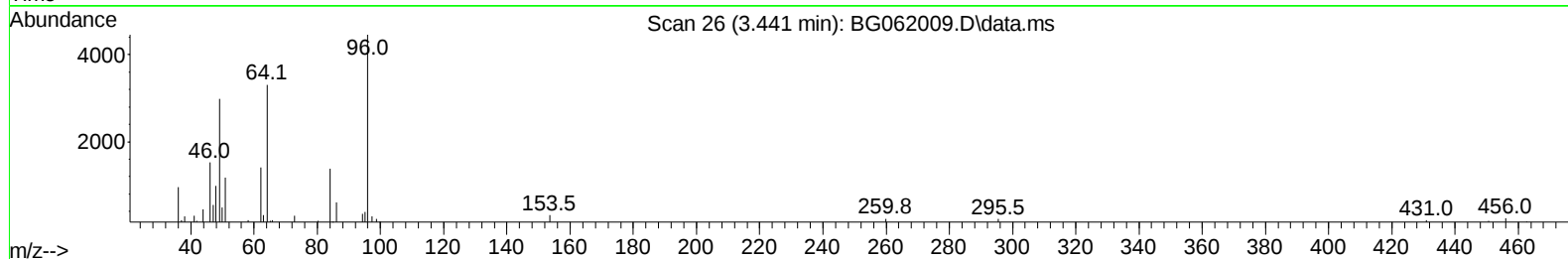
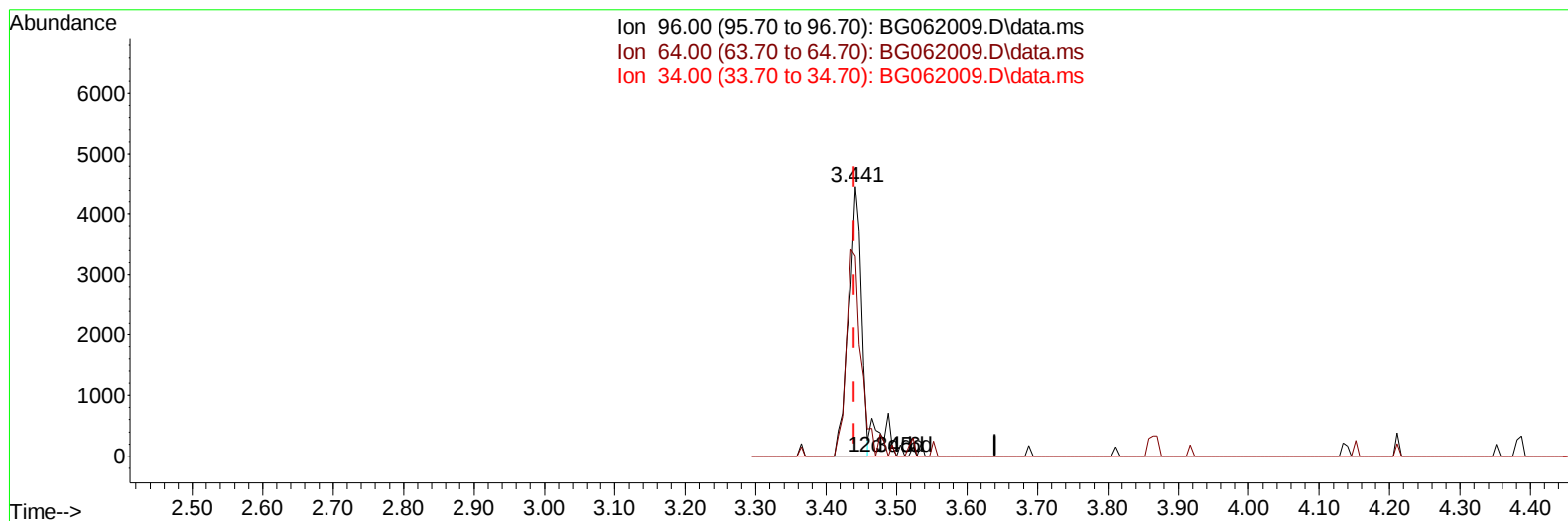
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071124\
Data File : BG062009.D
Acq On : 11 Jul 2024 14:13
Operator : MA/JU
Sample : PB161824BL
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
SBLK824

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 07/12/2024
Supervised By :mohammad ahmed 07/15/2024

Quant Time: Jul 12 00:06:02 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jul 11 12:11:32 2024
Response via : Initial Calibration



TIC: BG062009.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.441min (+ 0.001) 5.88 ng/uL

response 5563

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	104.80	74.16#
34.00	0.00	0.00
0.00	0.00	0.00

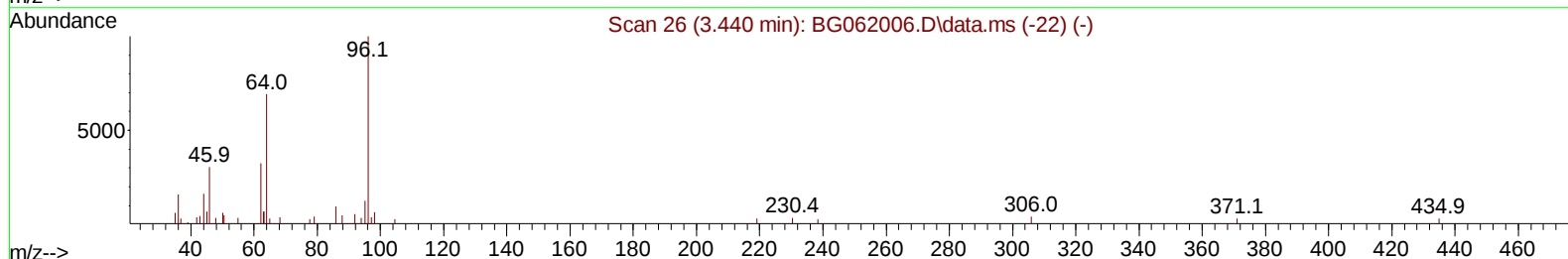
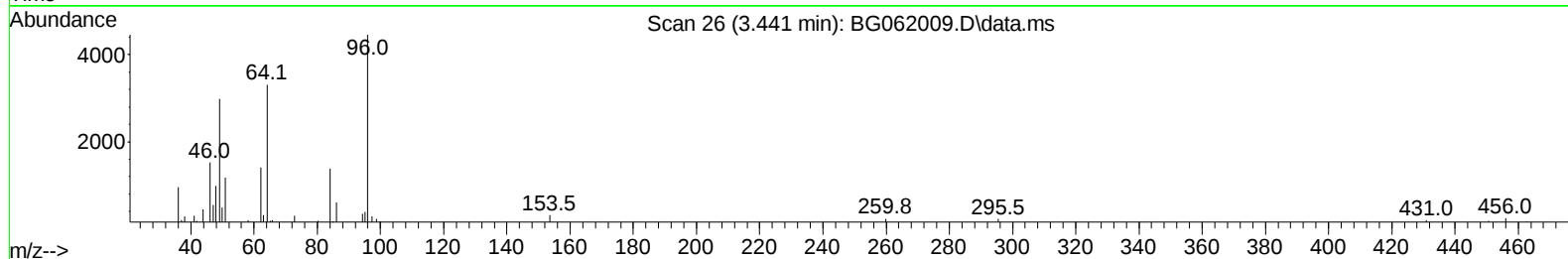
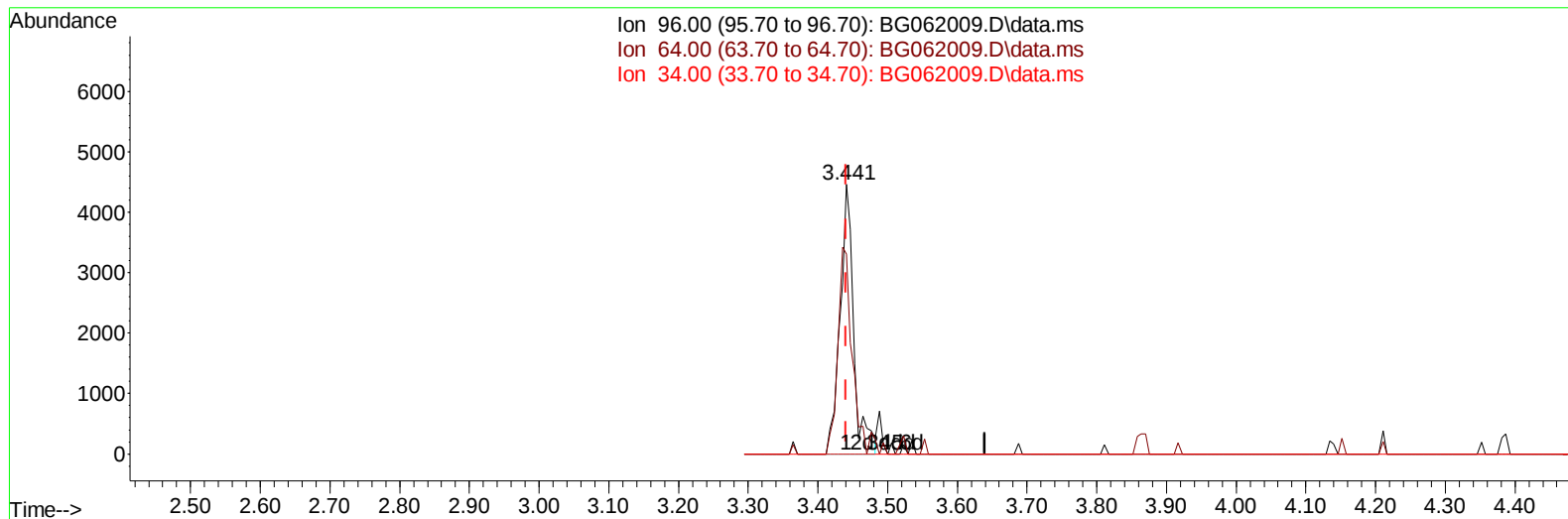
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TIC: BG062009.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.441min (+ 0.001) 6.51 ng/uL m

response 6157

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	104.80	74.16#
34.00	0.00	0.00
0.00	0.00	0.00

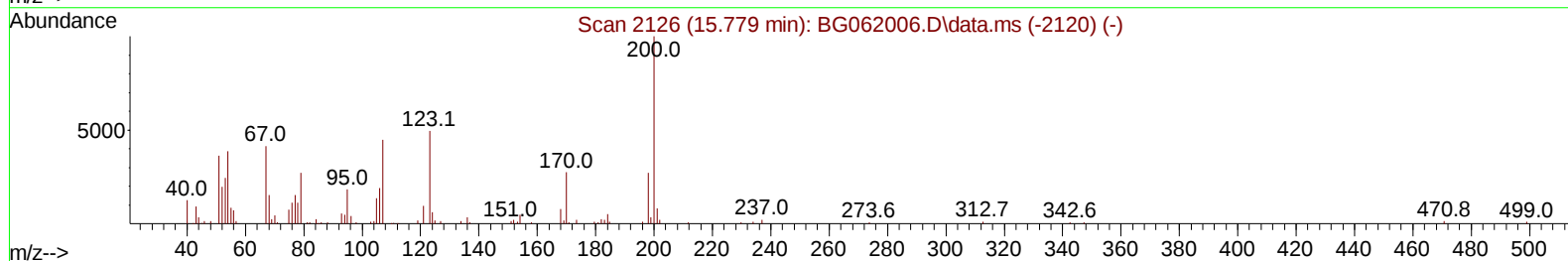
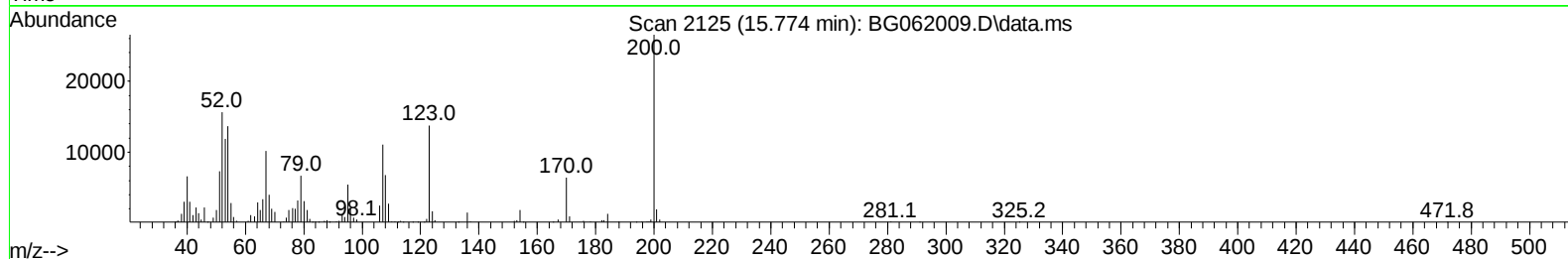
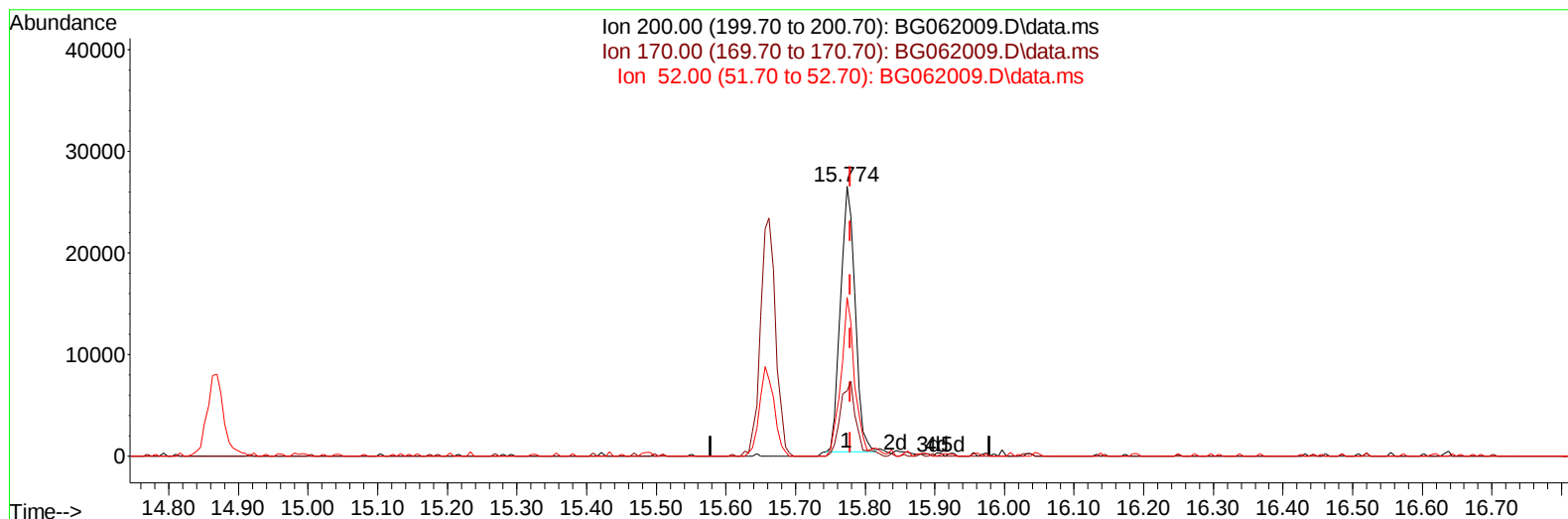
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 ALS Vial : 5 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

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TIC: BG062009.D\data.ms

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

15.774min (-0.005) 20.73 ng/ul

response 38456

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	21.10	24.19
52.00	66.90	58.94
0.00	0.00	0.00

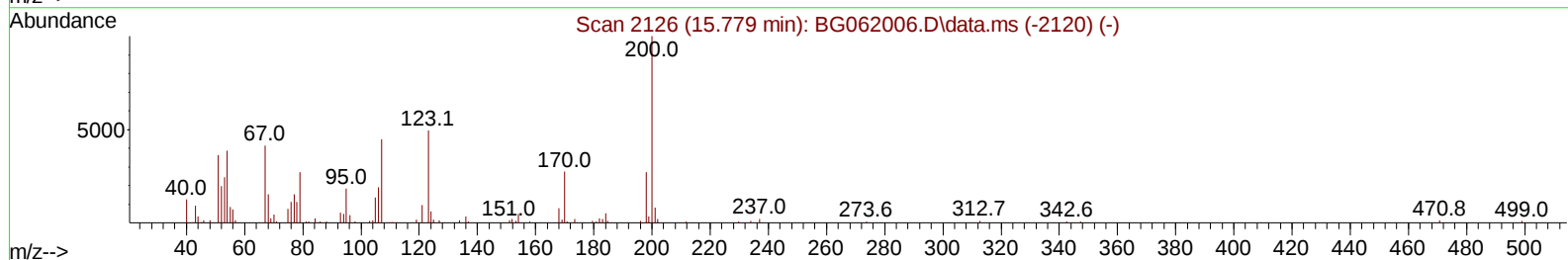
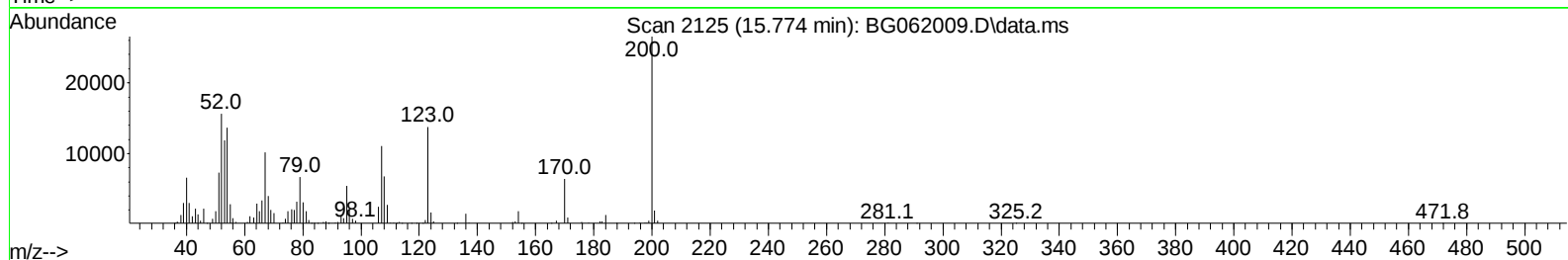
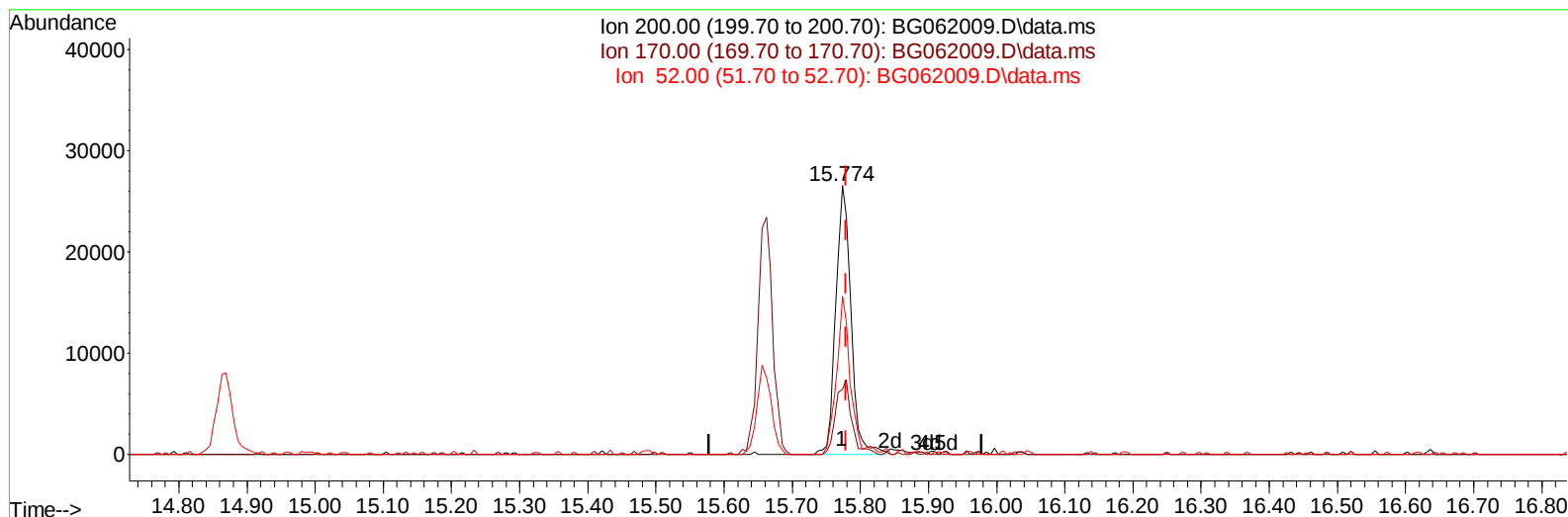
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Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Manual IntegrationsAPPROVED

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Supervised By :mohammad ahmed 07/15/2024

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TIC: BG062009.D\data.ms

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

15.774min (-0.005) 22.14 ng/ul m

response 41081

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	21.10	24.19
52.00	66.90	58.94
0.00	0.00	0.00

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 ALS Vial : 5 Sample Multiplier: 1

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 07/12/2024
 Supervised By :mohammad ahmed 07/15/2024

Quant Time: Jul 12 00:21:51 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.041	152	35558	20.000	ng/ul	0.00
20) Naphthalene-d8	10.850	136	187050	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.669	164	151557	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.413	188	353295	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.673	240	332815	20.000	ng/ul	0.00
88) Perylene-d12	24.875	264	373811	20.000	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.441	96	6157m	6.507	ng/uL	0.00
4) Pyridine-d5	3.858	84	71259	24.492	ng/ul	0.00
7) Phenol-d5	7.195	99	96788	24.873	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.360	67	61376	24.907	ng/ul	0.00
11) 2-Chlorophenol-d4	7.566	132	75459	28.270	ng/ul	0.00
15) 4-Methylphenol-d8	8.741	113	78179	25.517	ng/ul	0.00
21) Nitrobenzene-d5	9.205	128	35491	24.948	ng/ul	0.00
24) 2-Nitrophenol-d4	9.927	143	45140	27.786	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.468	165	89201	28.467	ng/ul	0.00
31) 4-Chloroaniline-d4	10.985	131	91082	20.021	ng/ul	0.00
46) Dimethylphthalate-d6	14.070	166	345810	27.754	ng/ul	0.00
49) Acenaphthylene-d8	14.363	160	351522	26.980	ng/ul	0.00
54) 4-Nitrophenol-d4	14.869	143	50708	25.484	ng/ul	0.00
60) Fluorene-d10	15.662	176	294233	28.051	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.774	200	41081m	22.141	ng/ul	0.00
73) Anthracene-d10	17.513	188	443887	26.911	ng/ul	0.00
81) Pyrene-d10	19.798	212	529416	27.044	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.663	264	540155	29.136	ng/ul	0.00

Target Compounds Qvalue

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

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