

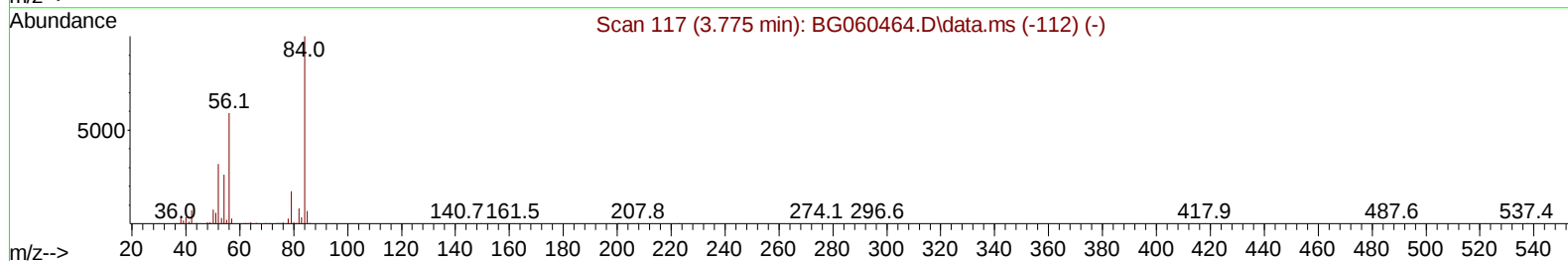
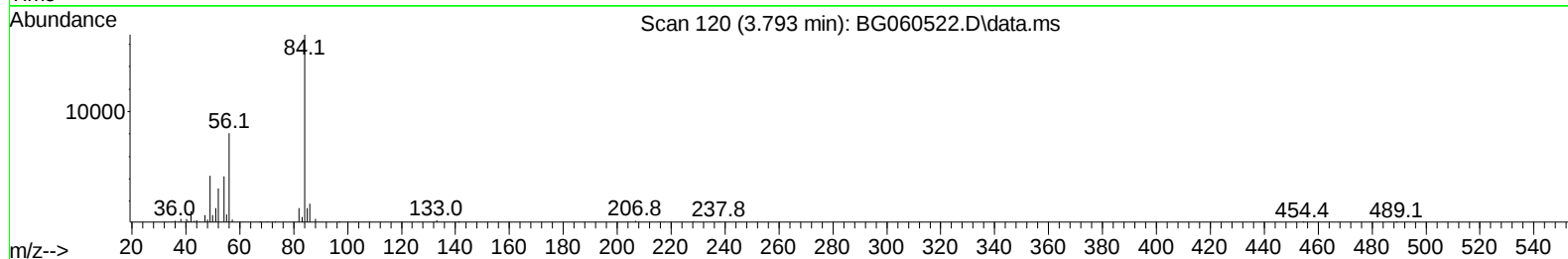
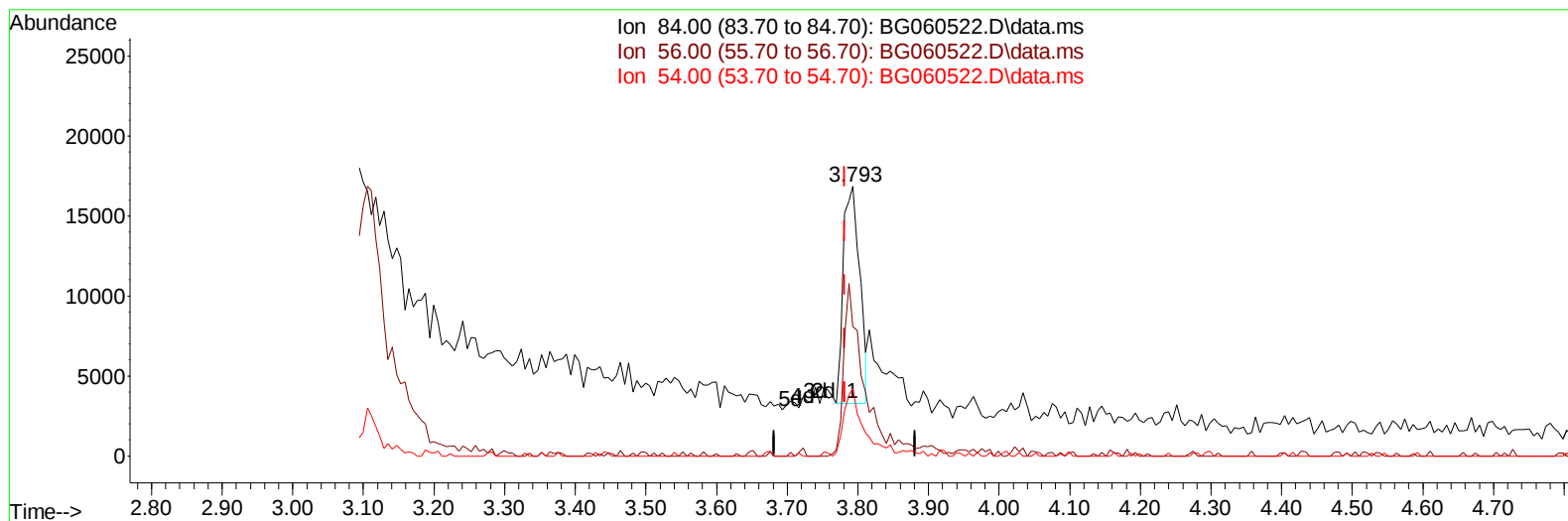
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG020124\
Data File : BG060522.D
Acq On : 31 Jan 2024 23:57
Operator : MA/JU
Sample : P1289-10
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
C0AF3

Manual IntegrationsAPPROVED

Quant Time: Feb 01 01:18:49 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG012624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jan 31 22:04:13 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/01/2024
Supervised By :Sohil Jodhani 02/02/2024



TIC: BG060522.D\data.ms

(4) Pyridine-d5 (S)

3.793min (+ 0.011) 2.94 ng/u1

response 21945

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	60.80	48.01#
54.00	26.10	25.15
0.00	0.00	0.00

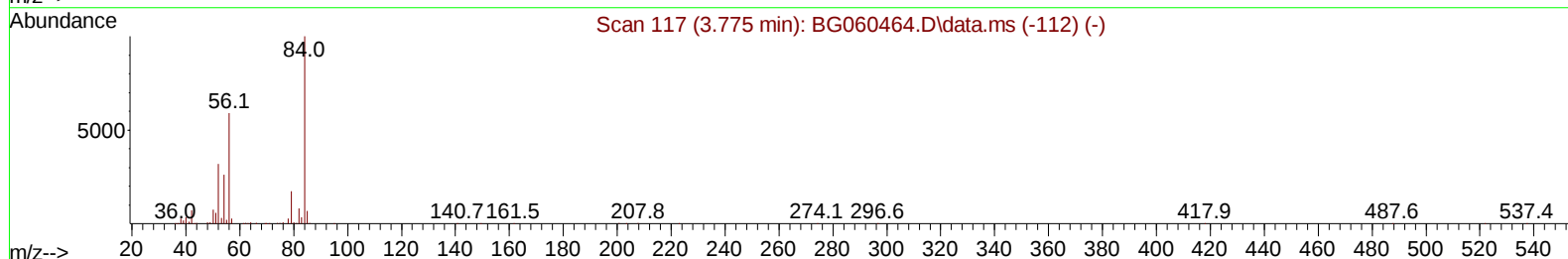
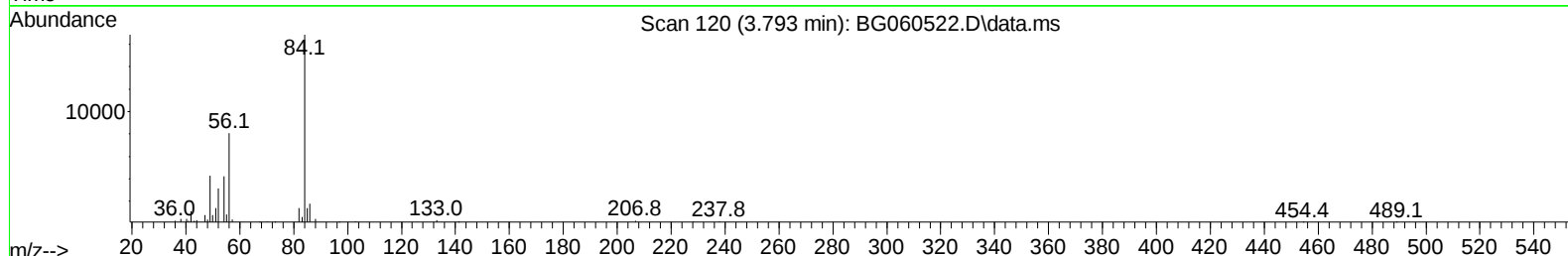
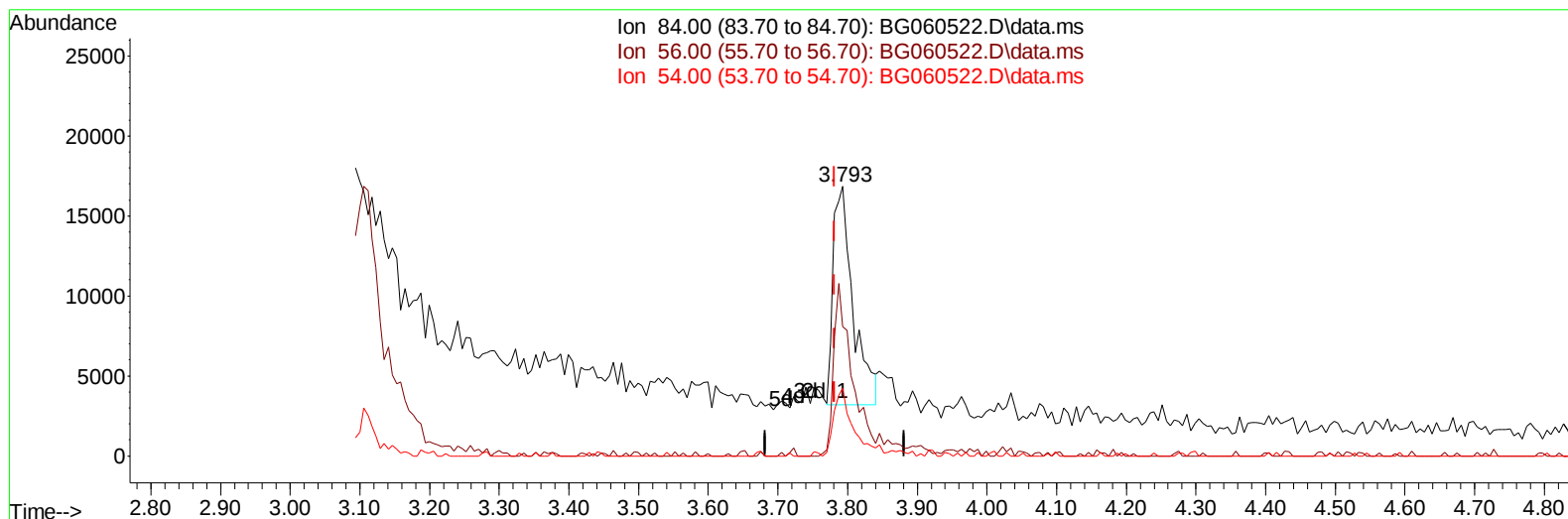
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TIC: BG060522.D\data.ms

(4) Pyridine-d5 (S)

3.793min (+ 0.011) 3.64 ng/ul m

response 27147

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	60.80	48.01#
54.00	26.10	25.15
0.00	0.00	0.00

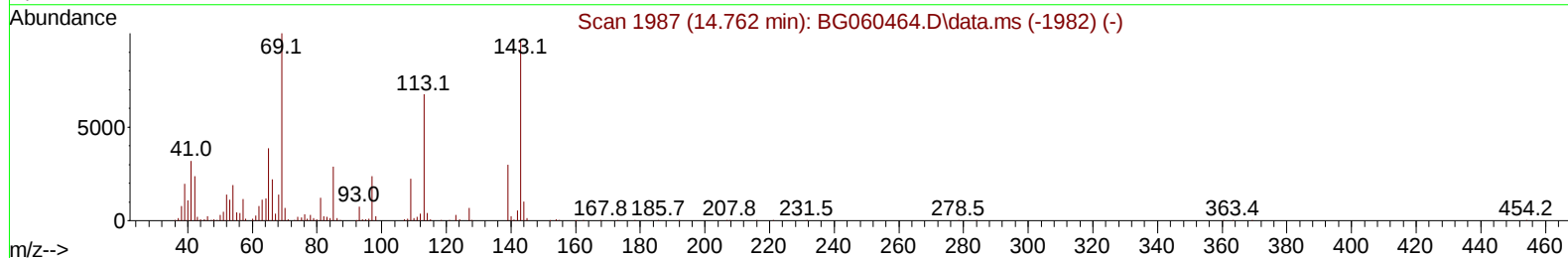
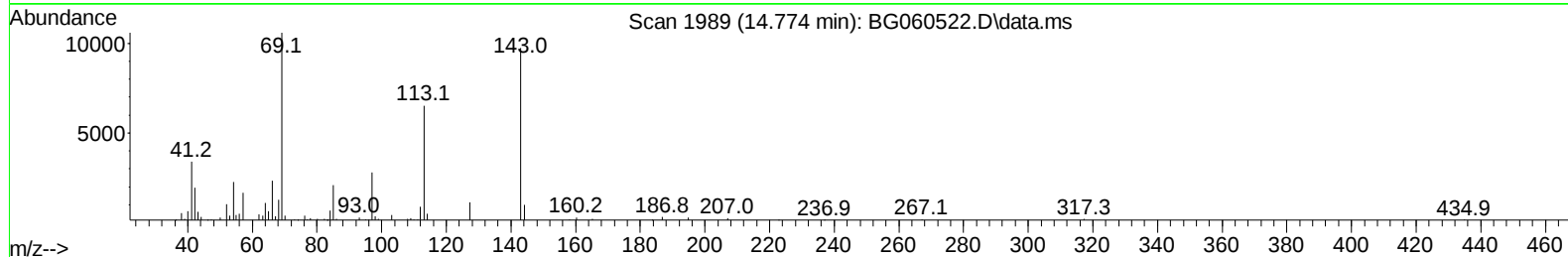
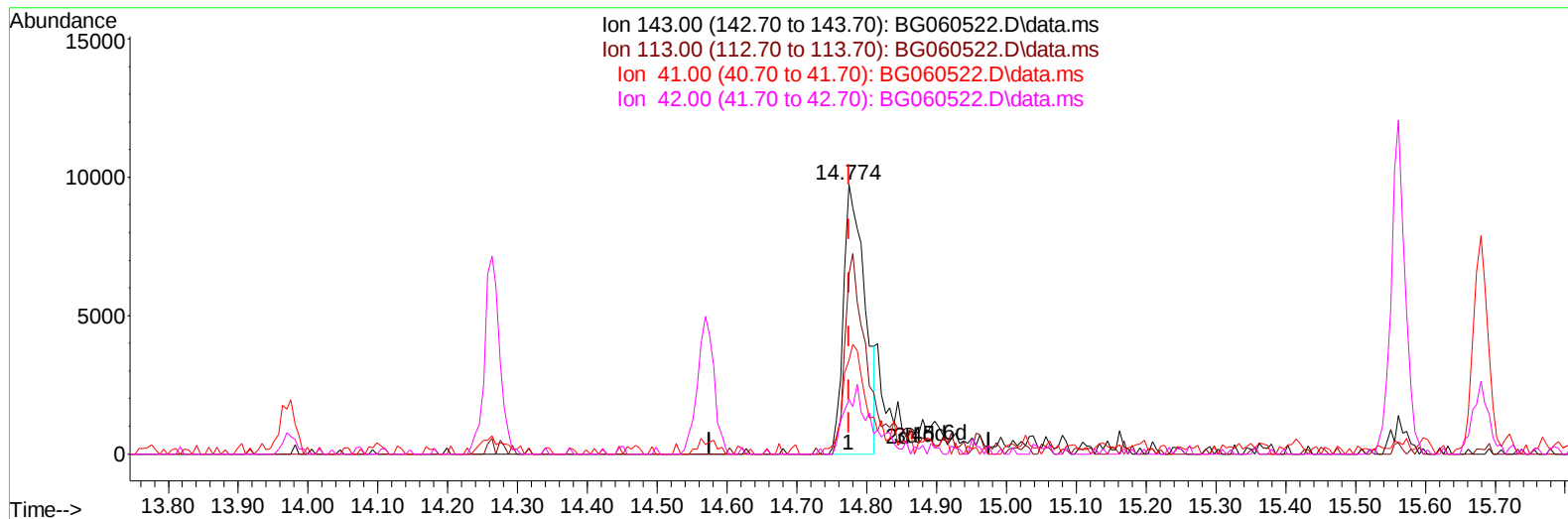
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG020124\
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Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: BG060522.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.774min (-0.001) 3.98 ng/u1

response 20525

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.60	66.94
41.00	29.80	34.84
42.00	22.70	20.22

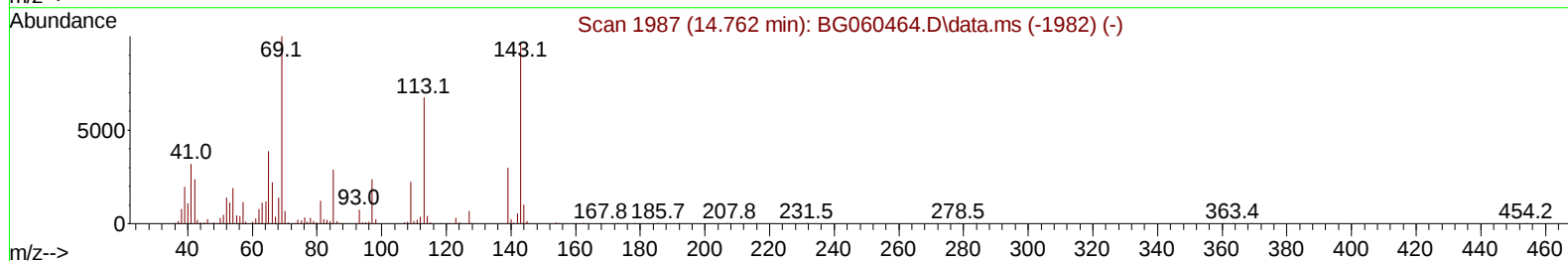
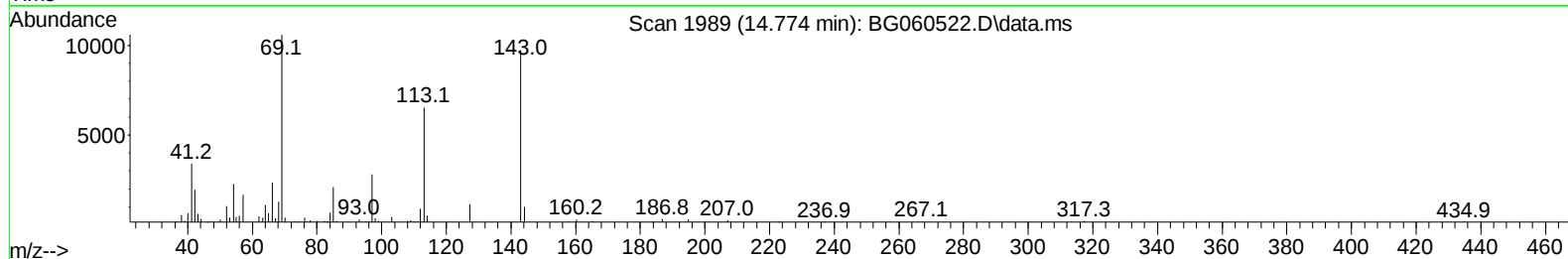
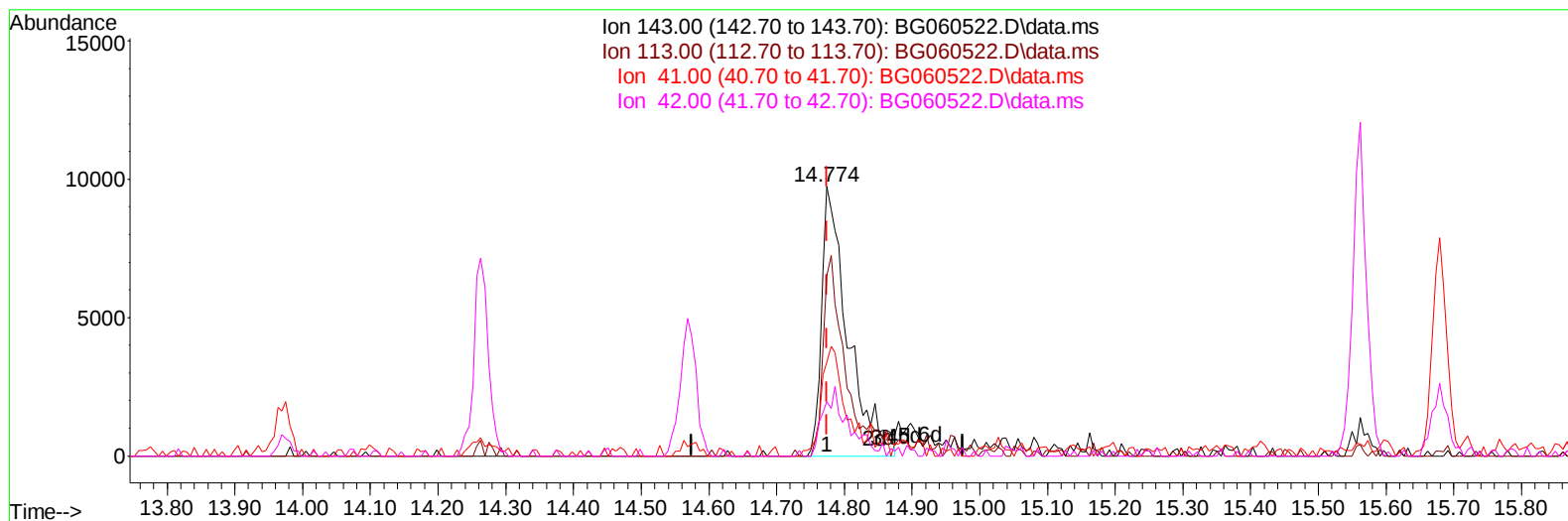
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Acq On : 31 Jan 2024 23:57
Operator : MA/JU
Sample : P1289-10
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BG060522.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.774min (-0.001) 5.12 ng/ul m

response 26367

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.60	66.94
41.00	29.80	34.84
42.00	22.70	20.22

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 ALS Vial : 13 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 02/01/2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.935	152	118155	20.000	ng/ul	0.00
20) Naphthalene-d8	10.738	136	516767	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.569	164	439497	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.318	188	1237689	20.000	ng/ul	0.00
79) Chrysene-d12	21.578	240	1190865	20.000	ng/ul	0.00
88) Perylene-d12	24.751	264	1327240	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.376	96	10734	4.552	ng/uL	0.00
4) Pyridine-d5	3.793	84	27147m	3.640	ng/ul	0.01
7) Phenol-d5	7.101	99	61547	5.485	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.260	67	203803	28.902	ng/ul	0.00
11) 2-Chlorophenol-d4	7.465	132	166361	22.200	ng/ul	0.00
15) 4-Methylphenol-d8	8.640	113	126891	13.728	ng/ul	0.00
21) Nitrobenzene-d5	9.099	128	132228	34.686	ng/ul	0.00
24) 2-Nitrophenol-d4	9.821	143	144877	32.529	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.362	165	275657	29.761	ng/ul	0.00
31) 4-Chloroaniline-d4	10.873	131	247662	20.382	ng/ul	0.00
46) Dimethylphthalate-d6	13.975	166	1272624	35.387	ng/ul	0.00
49) Acenaphthylene-d8	14.263	160	1395787	34.600	ng/ul	0.00
54) 4-Nitrophenol-d4	14.774	143	26367m	5.119	ng/ul	0.00
60) Fluorene-d10	15.562	176	1190181	37.770	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.679	200	217711	28.407	ng/ul	0.00
73) Anthracene-d10	17.418	188	2231595	38.036	ng/ul	0.00
81) Pyrene-d10	19.710	212	2758753	38.652	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.533	264	2763319	39.689	ng/ul	0.00

Target Compounds Qvalue

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

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