

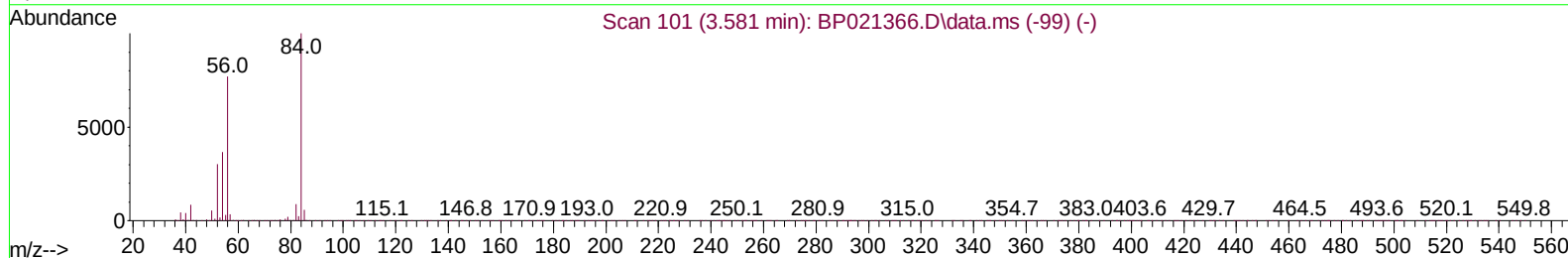
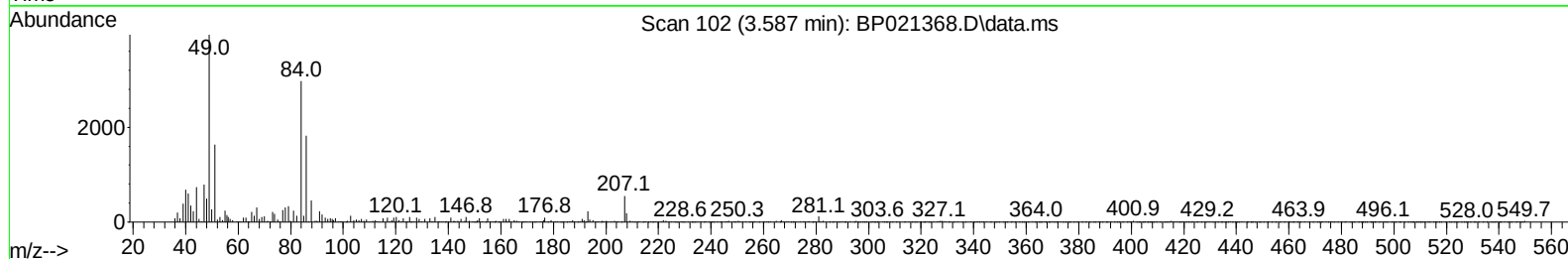
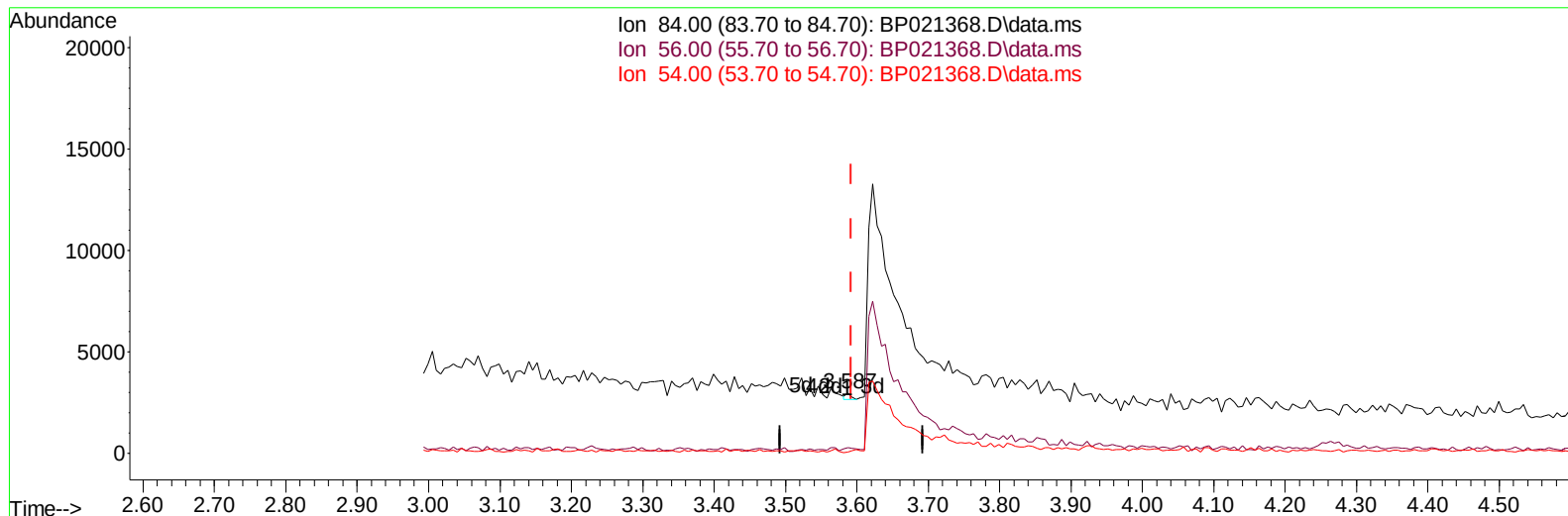
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080524\
Data File : BP021368.D
Acq On : 05 Aug 2024 11:31
Operator : CG/JU
Sample : P3300-04
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
E1GE7

Manual IntegrationsAPPROVED

Quant Time: Aug 05 12:00:59 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jul 29 12:46:18 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 08/06/2024
Supervised By :mohammad ahmed 08/07/2024



TIC: BP021368.D\data.ms

(4) Pyridine-d5 (S)

3.587min (-0.005) 0.02 ng/u1

response 156

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	70.20	8.72#
54.00	34.30	1.25#
0.00	0.00	0.00

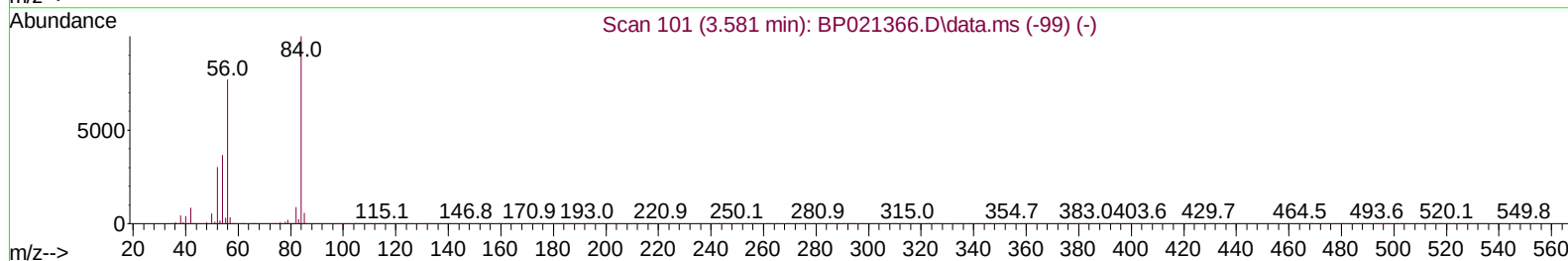
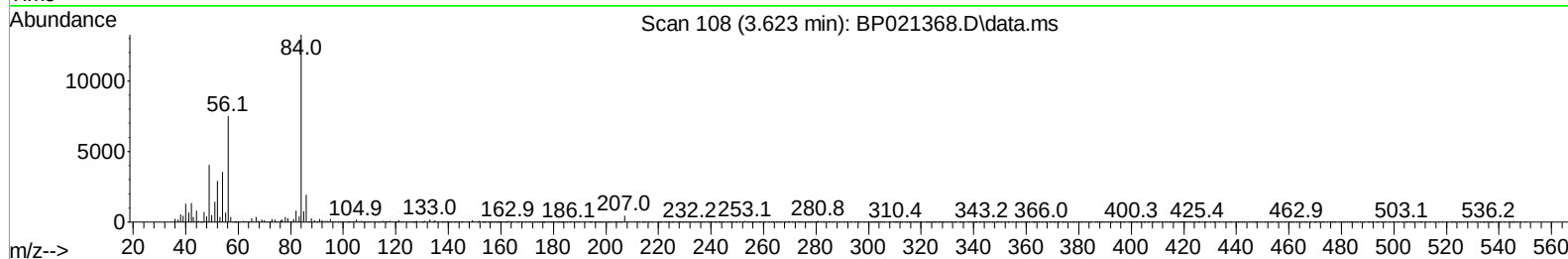
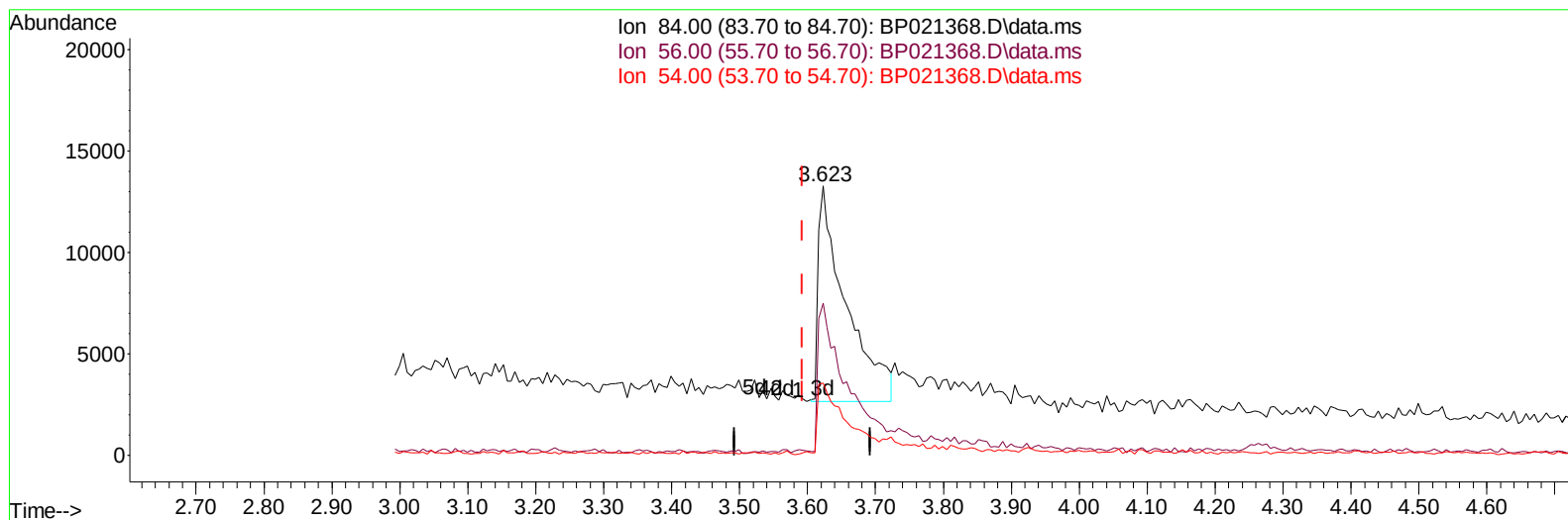
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(4) Pyridine-d5 (S)

3.623min (+ 0.030) 2.96 ng/u1 m

response 29875

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	70.20	56.55
54.00	34.30	26.67#
0.00	0.00	0.00

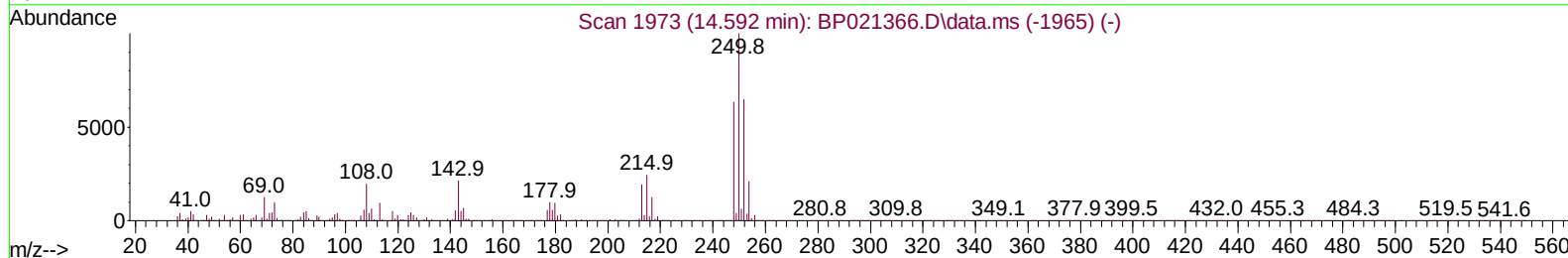
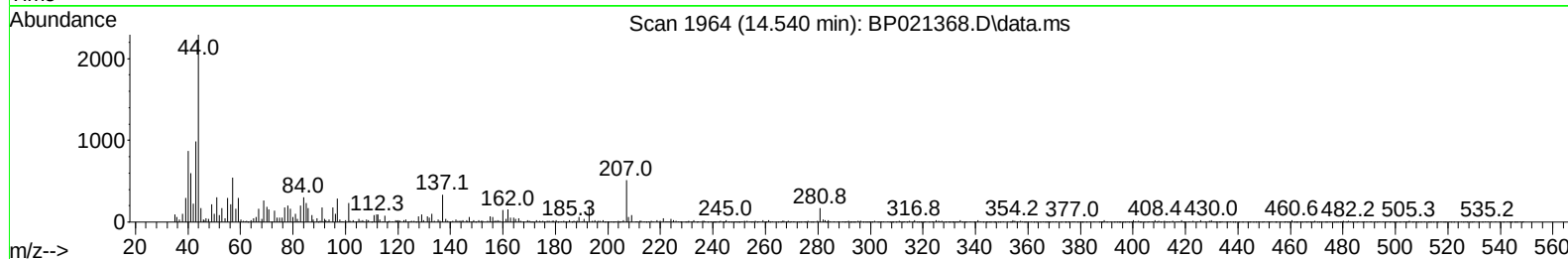
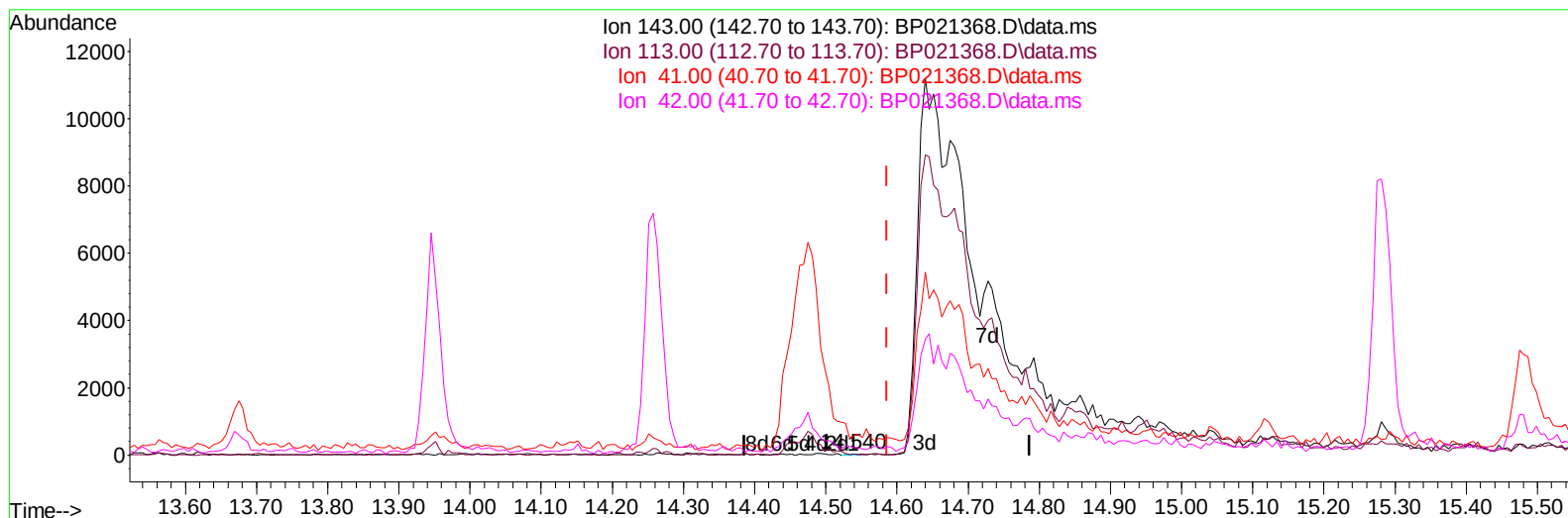
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 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BP021368.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.540min (-0.047) 0.01 ng/ul

response 26

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	87.88
41.00	37.90	1818.18#
42.00	27.10	675.76#

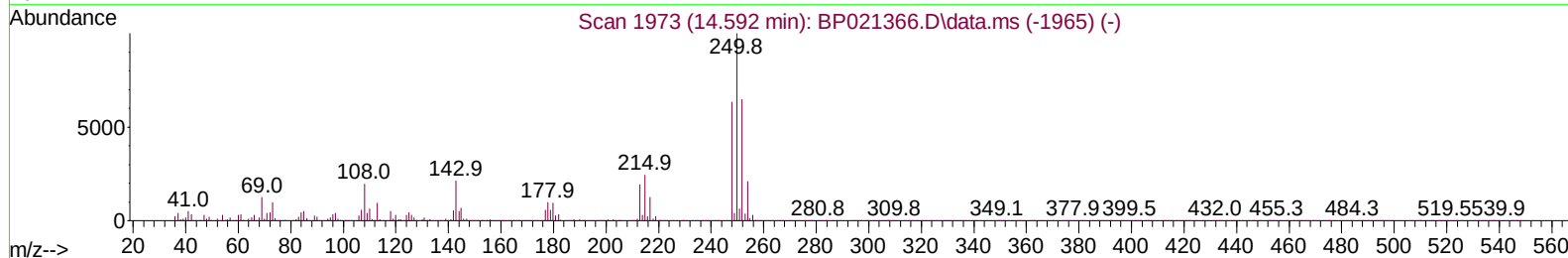
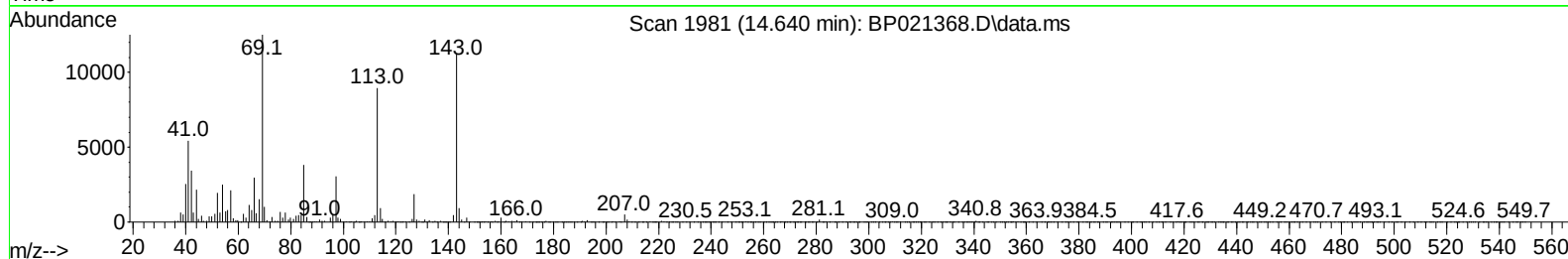
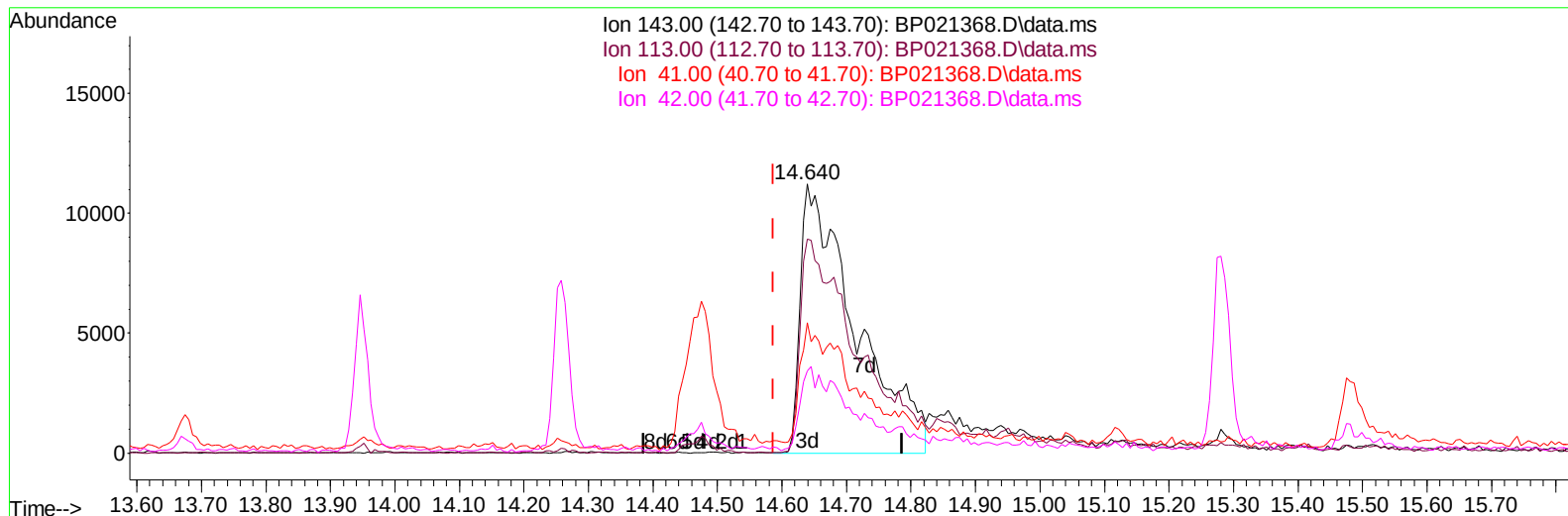
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Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
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TIC: BP021368.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.640min (+ 0.053) 16.42 ng/ul m

response 66562

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	74.70	79.64
41.00	37.90	48.46#
42.00	27.10	30.65

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 Sample : P3300-04
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E1GE7

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/06/2024
 Supervised By :mohammad ahmed 08/07/2024

Quant Time: Aug 05 12:02:02 2024
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 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jul 29 12:46:18 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.622	152	158702	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.381	136	623552	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.257	164	391905	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.081	188	872667	20.000	ng/ul	-0.01
79) Chrysene-d12	21.527	240	907210	20.000	ng/ul	0.00
88) Perylene-d12	24.821	264	1048531	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	11141	3.195	ng/uL	-0.01
4) Pyridine-d5	3.623	84	29875m	2.959	ng/ul	0.03
7) Phenol-d5	6.852	99	251046	19.403	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.975	67	179980	22.953	ng/ul	0.00
11) 2-Chlorophenol-d4	7.175	132	221034	20.734	ng/ul	0.00
15) 4-Methylphenol-d8	8.375	113	211452	19.677	ng/ul	0.01
21) Nitrobenzene-d5	8.793	128	110041	22.721	ng/ul	0.00
24) 2-Nitrophenol-d4	9.510	143	108501	19.573	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.069	165	167935	16.754	ng/ul	0.02
31) 4-Chloroaniline-d4	10.569	131	269774	20.316	ng/ul	0.01
46) Dimethylphthalate-d6	13.681	166	659540	23.150	ng/ul	0.00
49) Acenaphthylene-d8	13.946	160	740690	23.154	ng/ul	-0.02
54) 4-Nitrophenol-d4	14.640	143	66562m	16.421	ng/ul	0.05
60) Fluorene-d10	15.281	176	577044	24.689	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.481	200	65364	12.379	ng/ul	0.02
73) Anthracene-d10	17.187	188	952951	24.586	ng/ul	0.00
81) Pyrene-d10	19.575	212	1209294	21.561	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.586	264	1323548	25.594	ng/ul	-0.01

Target Compounds Qvalue

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

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