

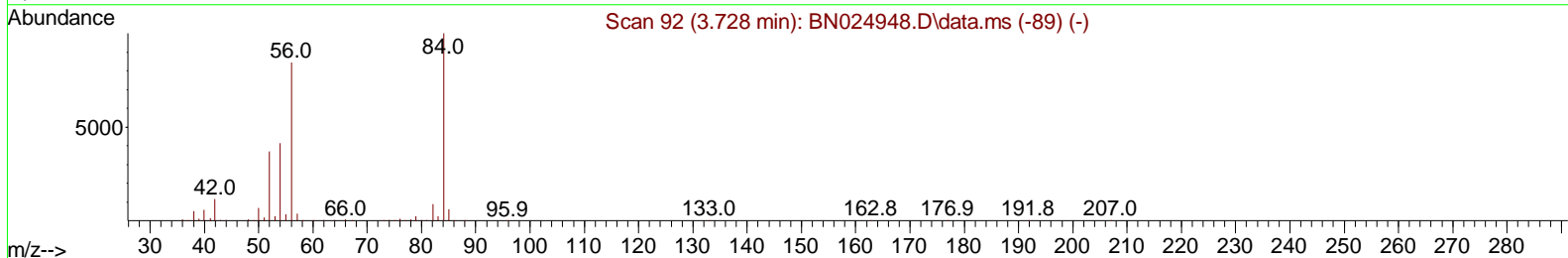
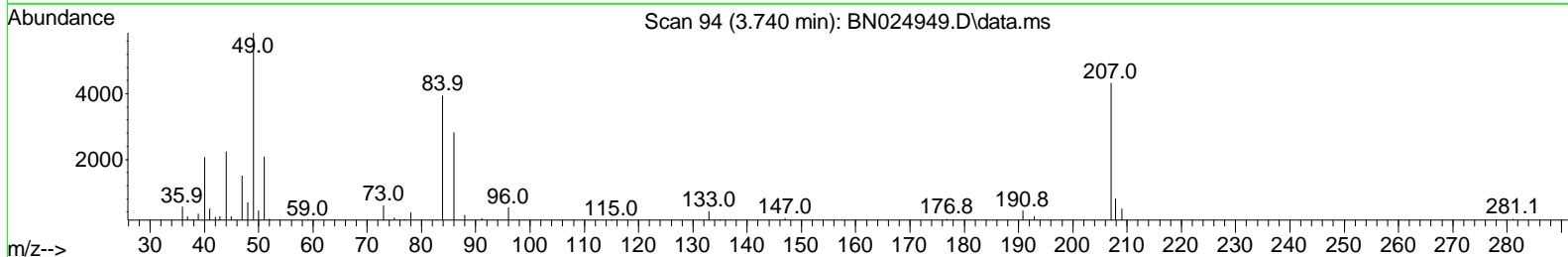
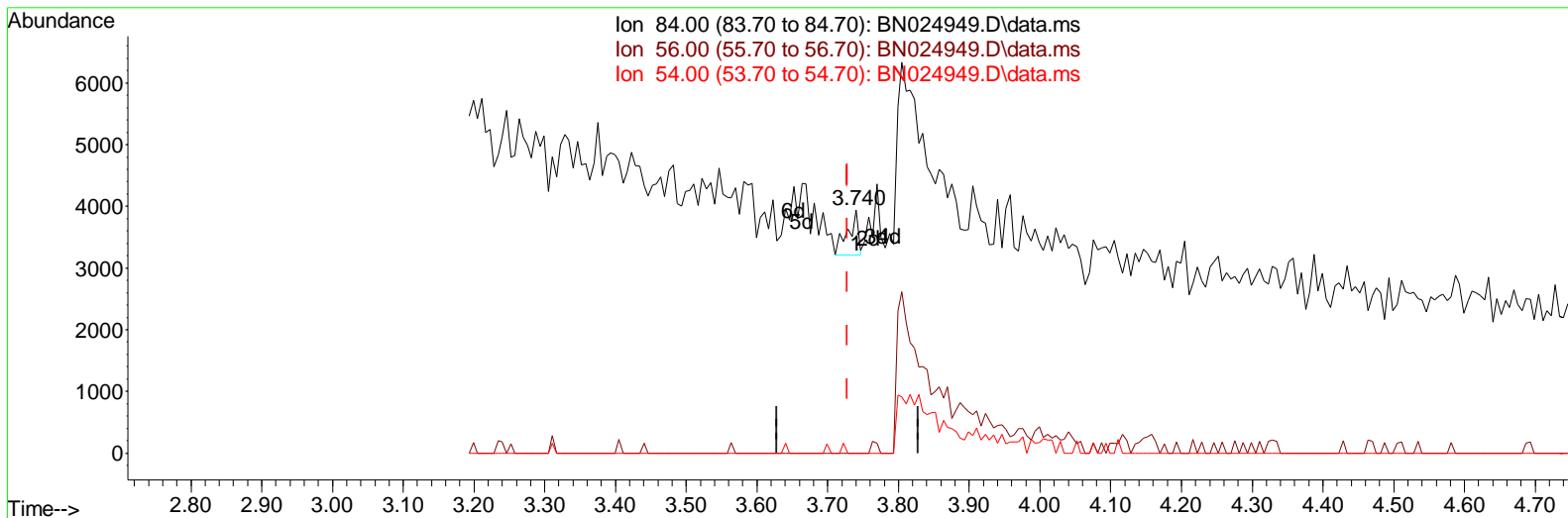
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041723\
 Data File : BN024949.D
 Acq On : 17 Apr 2023 11:07
 Operator : CG/JU
 Sample : PB152129BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
ClientSampleId :
 SBLK129

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 04/18/2023
 Supervised By :Jagrut Upadhyay 04/18/2023

Quant Time: Apr 18 01:30:50 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041023.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Apr 18 01:19:27 2023
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TIC: BN024949.D\data.ms

(4) Pyridine-d5 (s)

3.740min (+ 0.012) 0.02 ng/ul

response 735

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	82.10	0.00#
54.00	40.20	0.00#
0.00	0.00	0.00

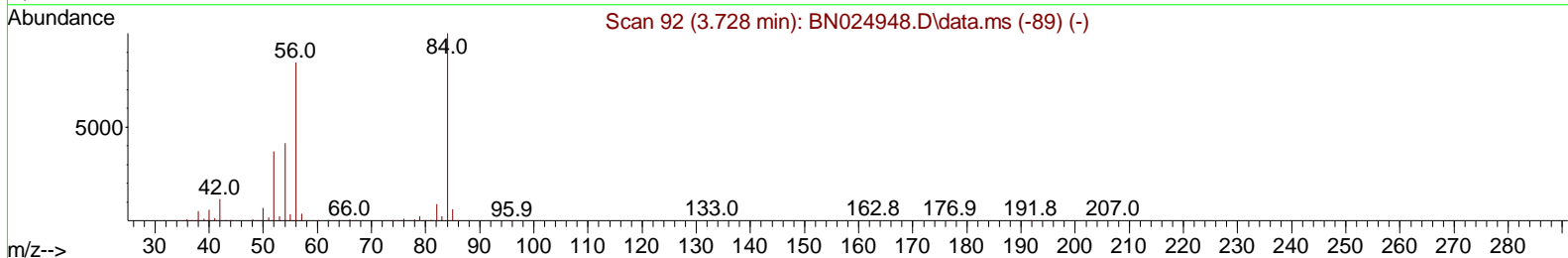
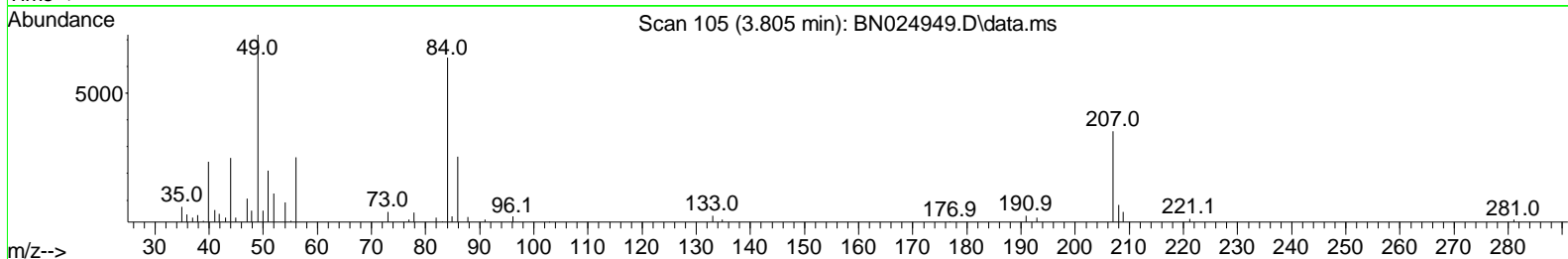
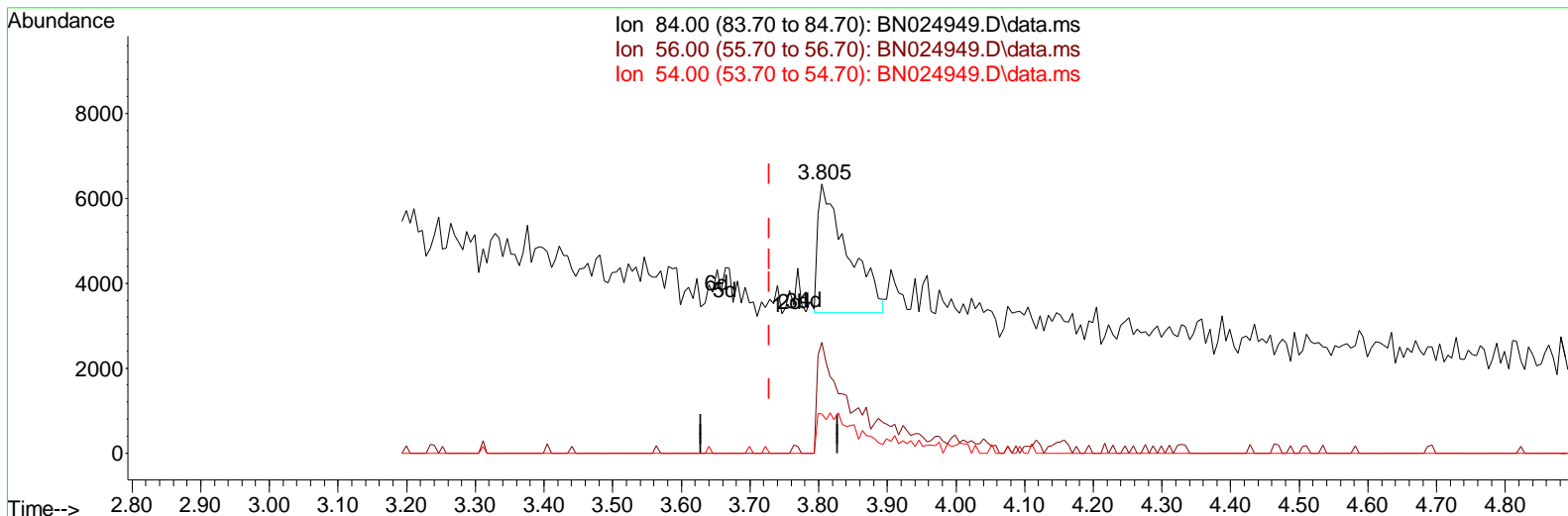
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(4) Pyridine-d5 (S)

3.805min (+ 0.077) 0.30 ng/ul m

response 9136

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	82.10	41.25#
54.00	40.20	14.45#
0.00	0.00	0.00

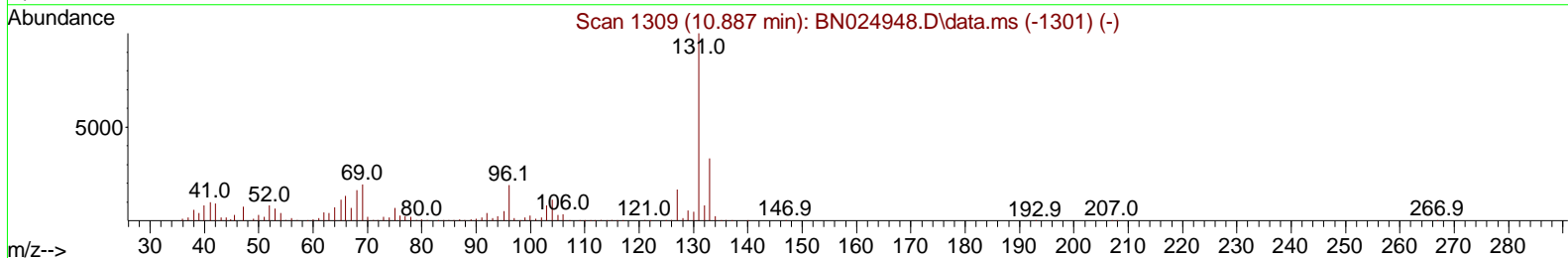
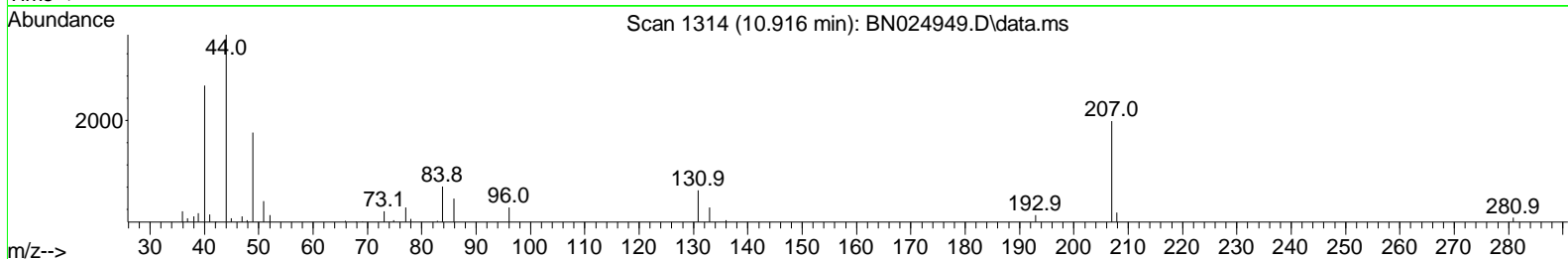
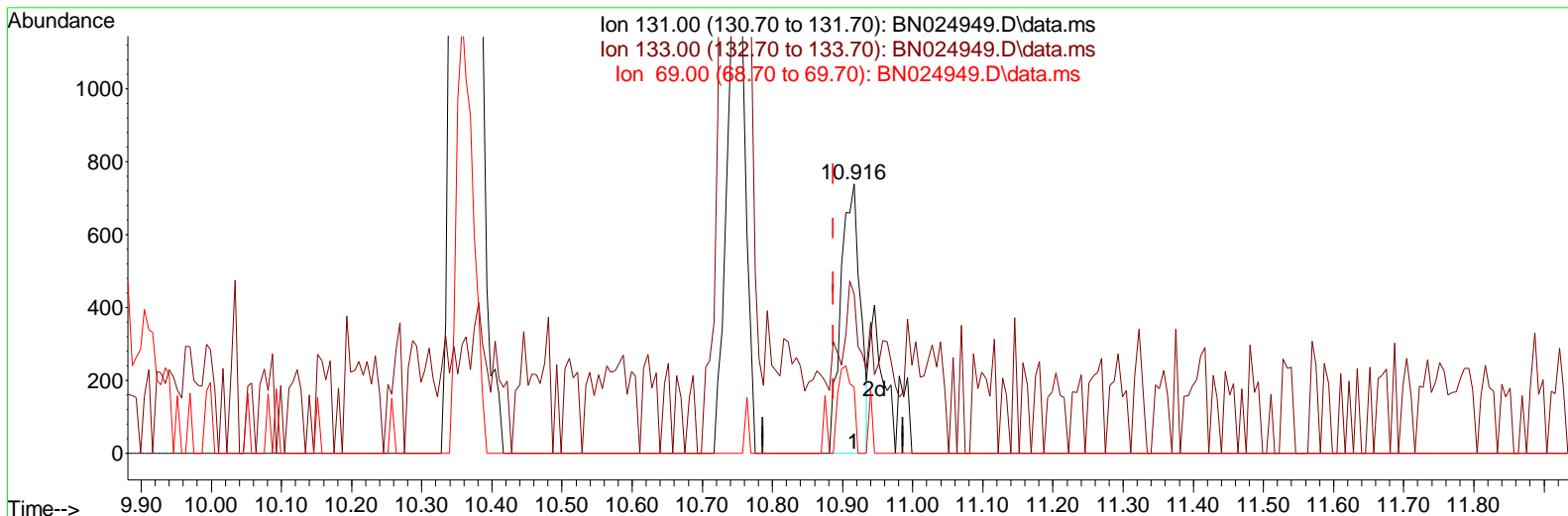
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 Sample : PB152129BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
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TIC: BN024949.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.916min (+ 0.030) 0.04 ng/ul

response 1425

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.00	59.00#
69.00	19.40	24.36#
0.00	0.00	0.00

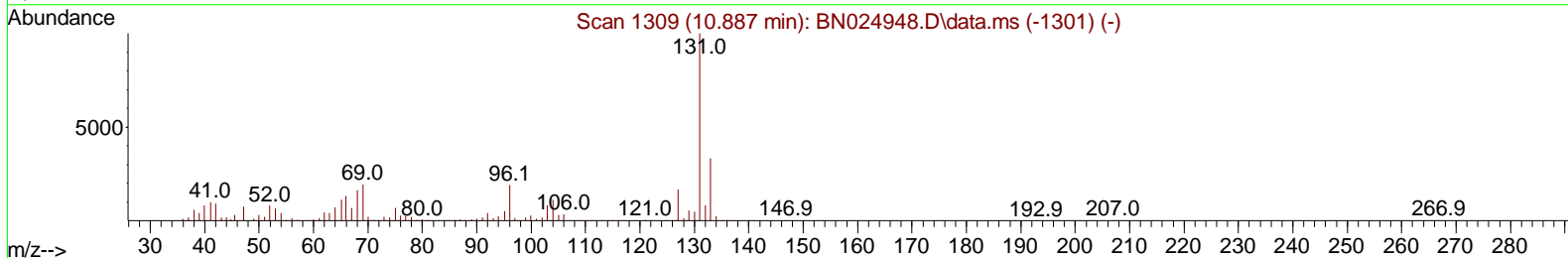
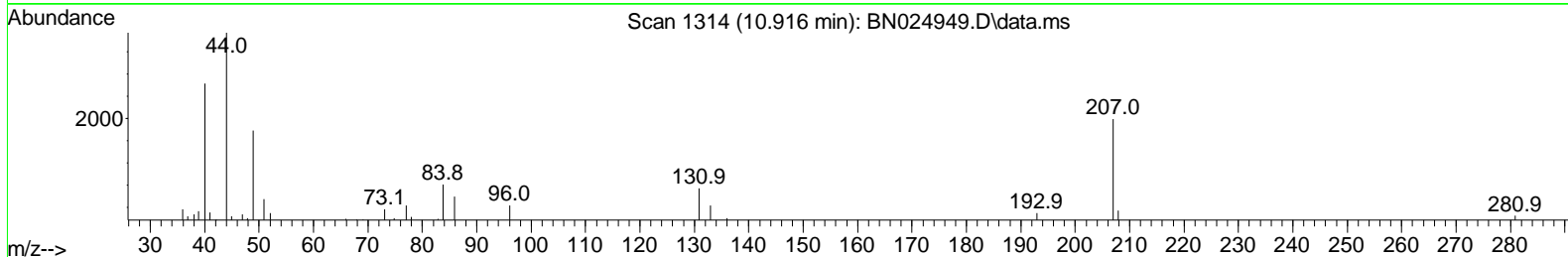
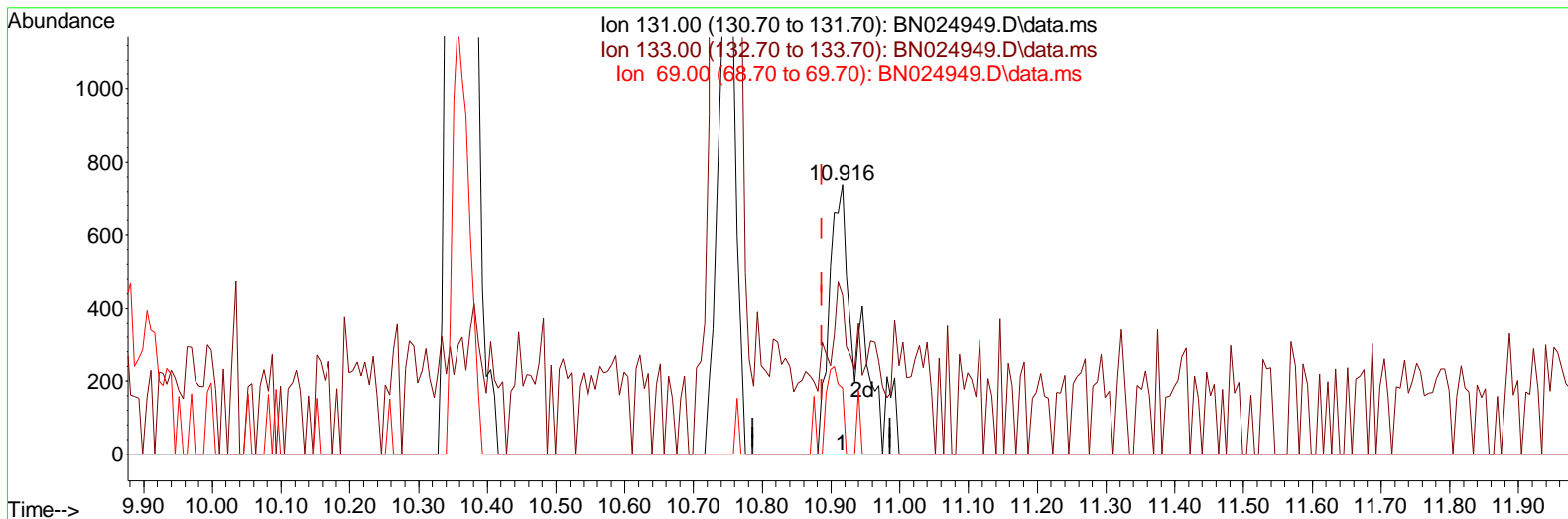
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 Sample : PB152129BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
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TIC: BN024949.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.916min (+ 0.030) 0.05 ng/ul m

response 1966

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.00	59.00#
69.00	19.40	24.36#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BNO41723\
 Data File : BNO24949.D
 Acq On : 17 Apr 2023 11:07
 Operator : CG/JU
 Sample : PB152129BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
ClientSampleId :
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Manual Integrations APPROVED

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 Quant Title : SVOA CALIBRATION
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 Response via : Initial Calibration

Compound	R. T.	QI on	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	386965	20.000	ng/ul	0.00
20) Naphthalene-d8	10.746	136	1675201	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.575	164	1049771	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.322	188	2294017	20.000	ng/ul	0.00
79) Chrysene-d12	21.527	240	2170682	20.000	ng/ul	0.00
88) Perylene-d12	23.963	264	2264451	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	48828	4.802	ng/uL	0.00
4) Pyridine-d5	3.805	84	9136m	0.300	ng/ul	0.08
7) Phenol-d5	7.093	99	823385	21.825	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.258	67	596592	23.643	ng/ul	0.00
11) 2-Chlorophenol-d4	7.458	132	659624	23.792	ng/ul	0.00
15) 4-Methylphenol-d8	8.640	113	648416	21.650	ng/ul	0.00
21) Nitrobenzene-d5	9.099	128	327234	24.125	ng/ul	0.00
24) 2-Nitrophenol-d4	9.822	143	358677	24.413	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	606783	23.107	ng/ul	0.00
31) 4-Chloroaniline-d4	10.916	131	1966m	0.049	ng/ul	0.03
46) Dimethylphthalate-d6	13.987	166	2044842	25.127	ng/ul	0.00
49) Acenaphthylene-d8	14.269	160	2235294	23.969	ng/ul	0.00
54) 4-Nitrophenol-d4	14.775	143	326078	19.547	ng/ul	0.00
60) Fluorene-d10	15.569	176	1694655	24.664	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.687	200	297431	19.923	ng/ul	0.00
73) Anthracene-d10	17.422	188	2755468	25.575	ng/ul	0.00
81) Pyrene-d10	19.722	212	3201307	25.398	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.810	264	2963106	25.026	ng/ul	-0.01

Target Compounds Qvalue

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

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