

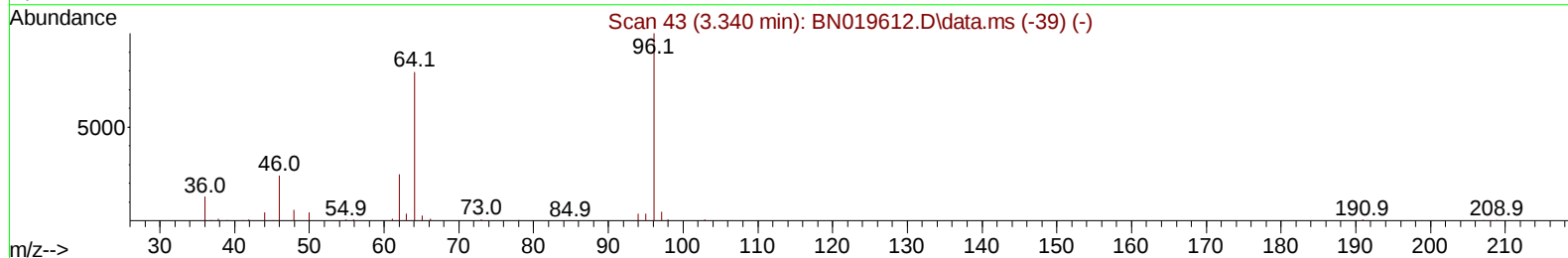
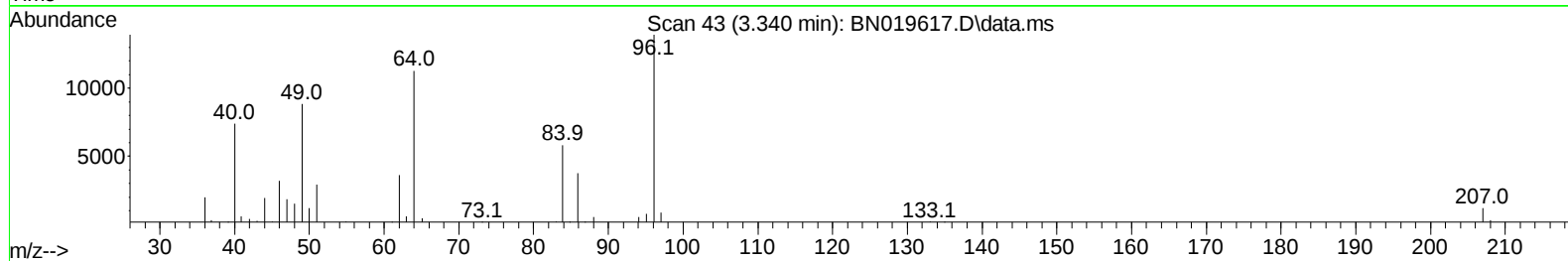
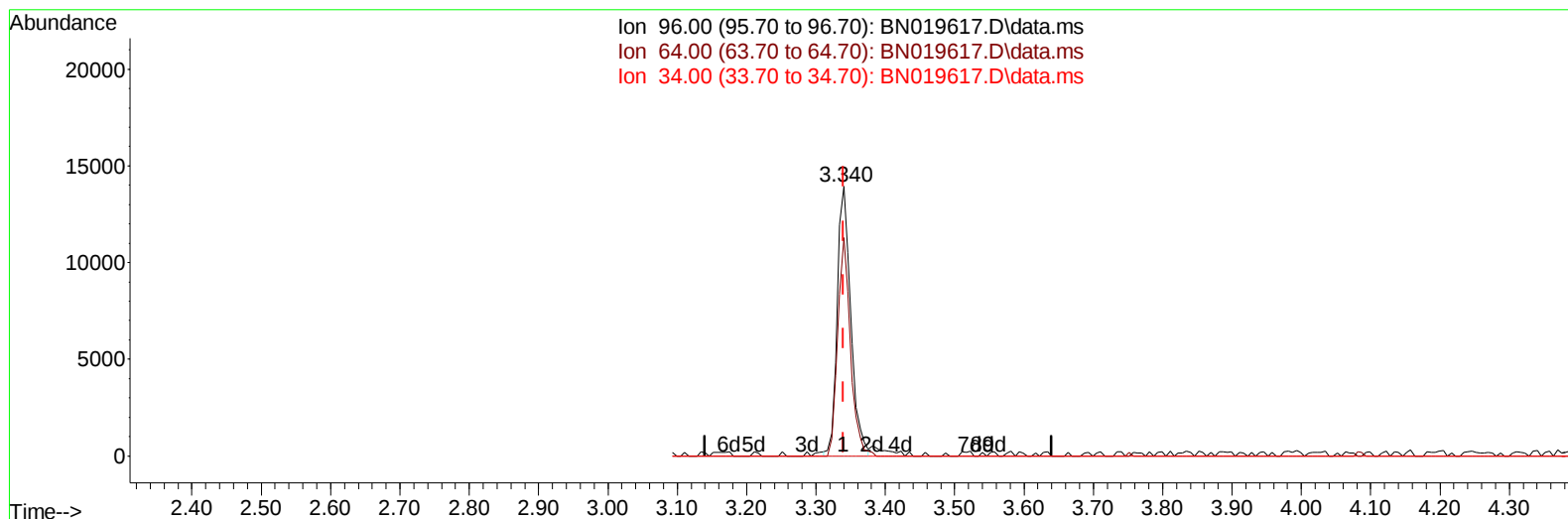
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN050222\
 Data File : BN019617.D
 Acq On : 02 May 2022 13:48
 Operator : CG/JU
 Sample : N2621-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BFFH9

Manual IntegrationsAPPROVED

Quant Time: May 03 03:46:55 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN042722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 03 03:31:44 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/05/2022
 Supervised By :Yogesh Patel 05/09/2022



TIC: BN019617.D\data.ms

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

3.340min (+ 0.000) 4.58 ng/uL

response 18884

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	80.20	80.91
34.00	0.00	0.00
0.00	0.00	0.00

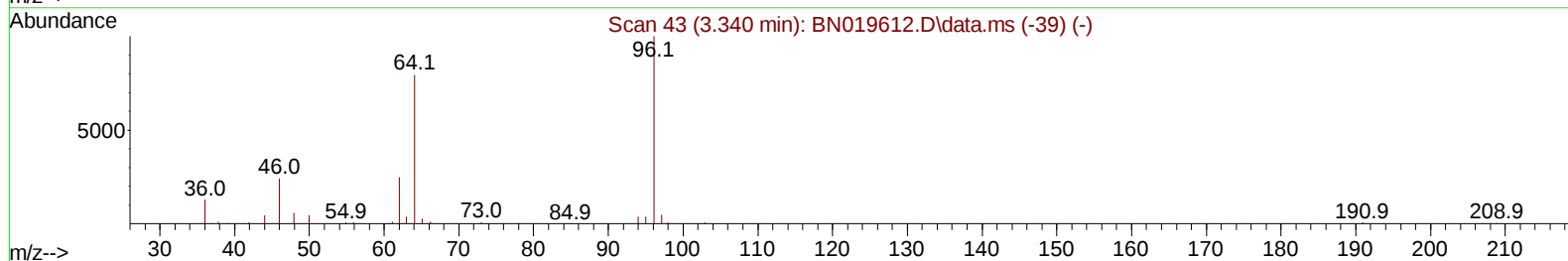
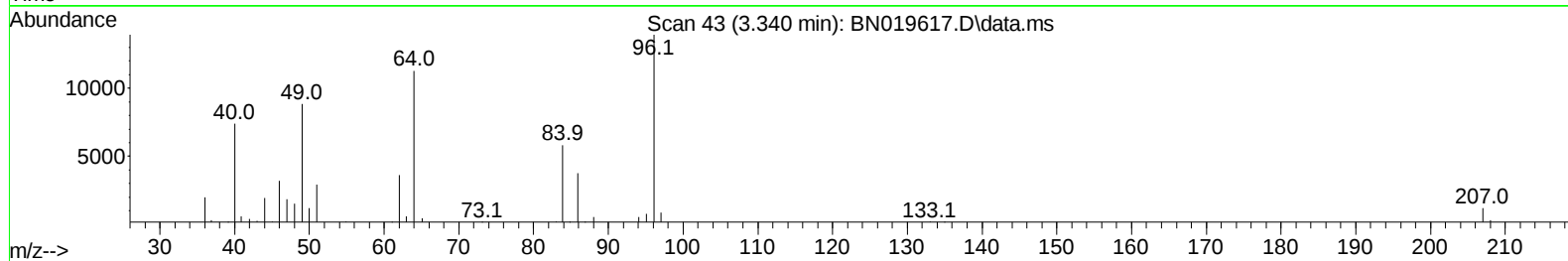
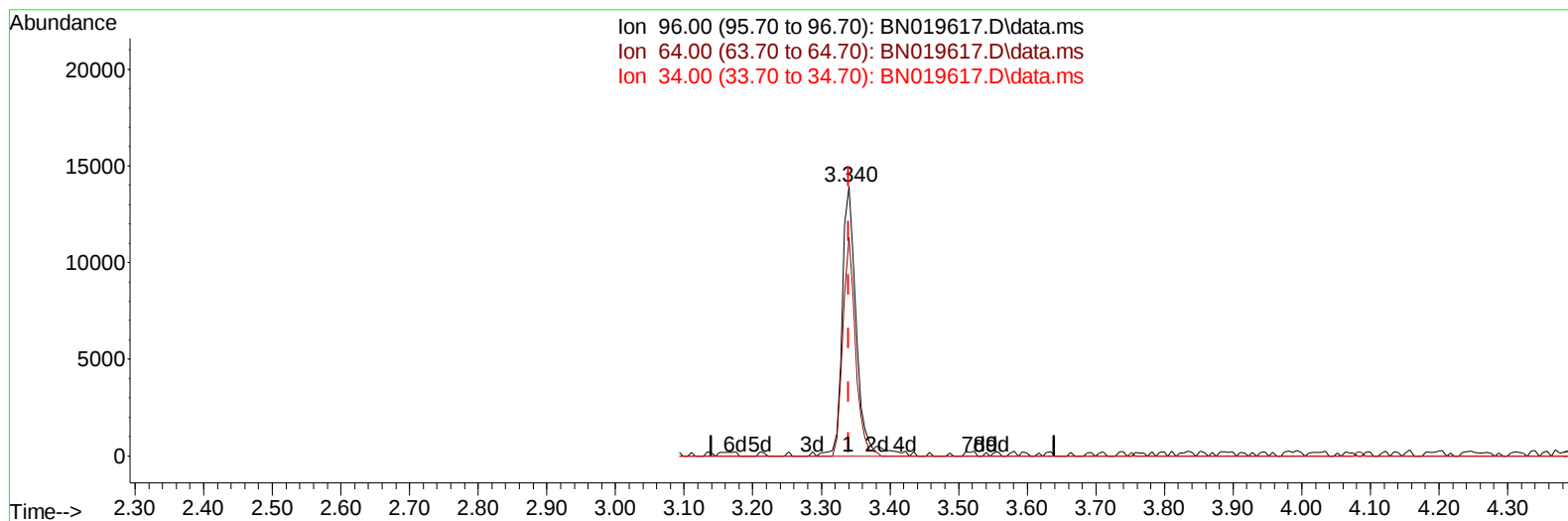
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TIC: BN019617.D\data.ms

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

3.340min (+ 0.000) 4.72 ng/uL m

response 19497

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	80.20	80.91
34.00	0.00	0.00
0.00	0.00	0.00

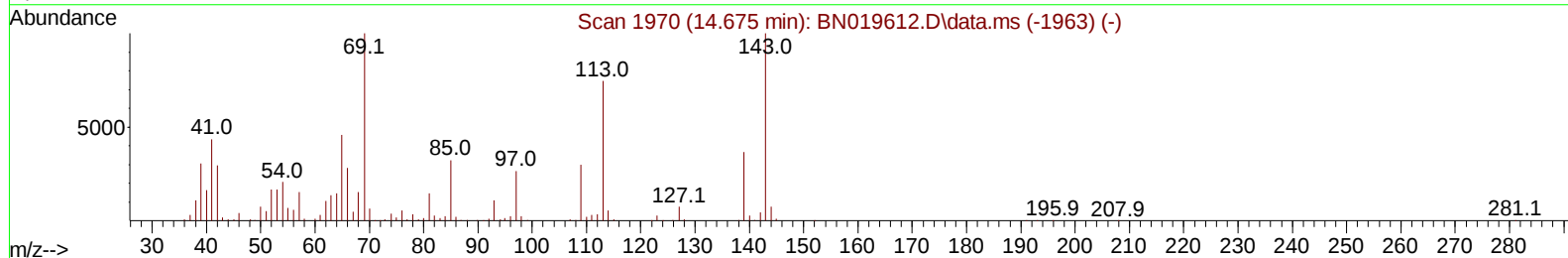
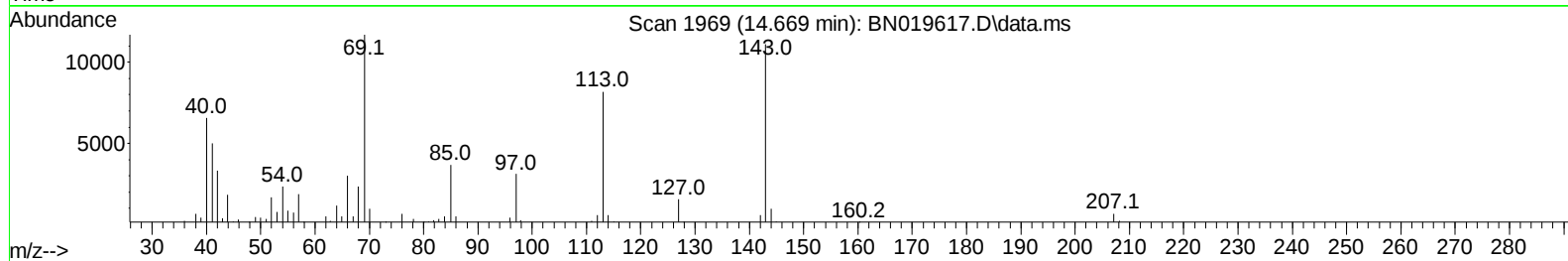
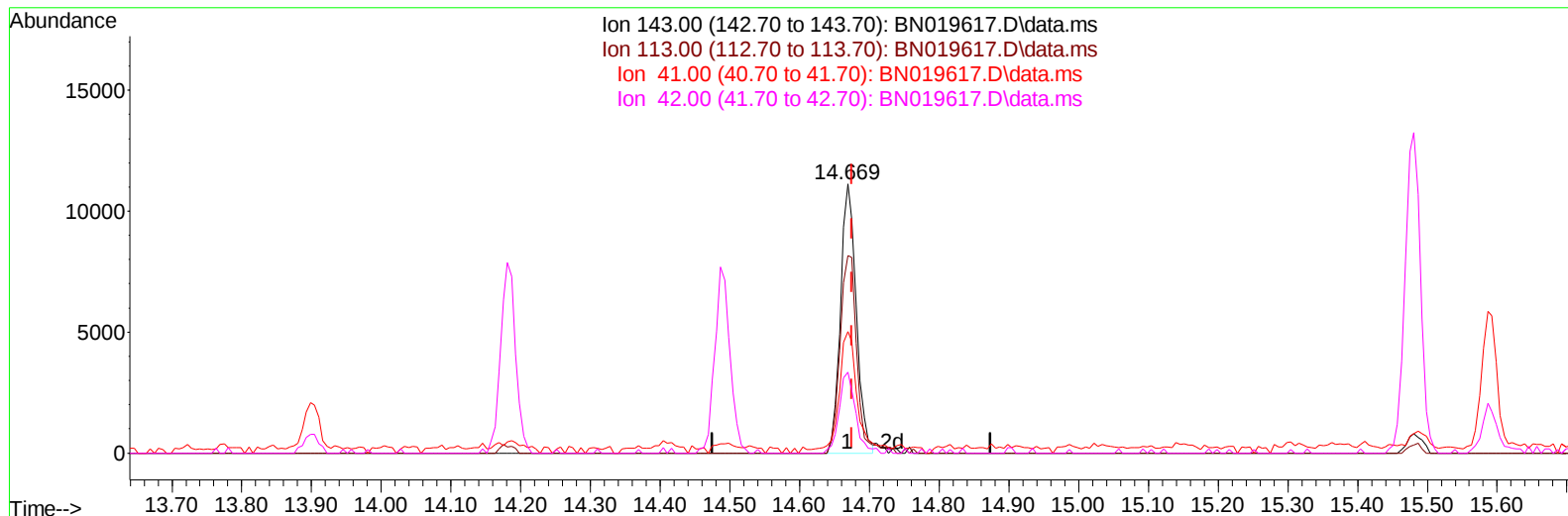
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ALS Vial : 7 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: BN019617.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (-0.006) 3.66 ng/u1

response 17302

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.20	73.55
41.00	39.70	45.09
42.00	27.00	30.04

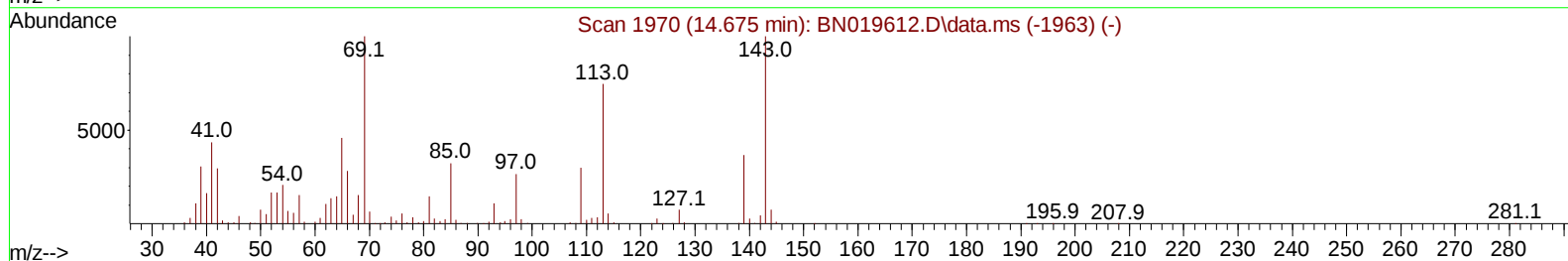
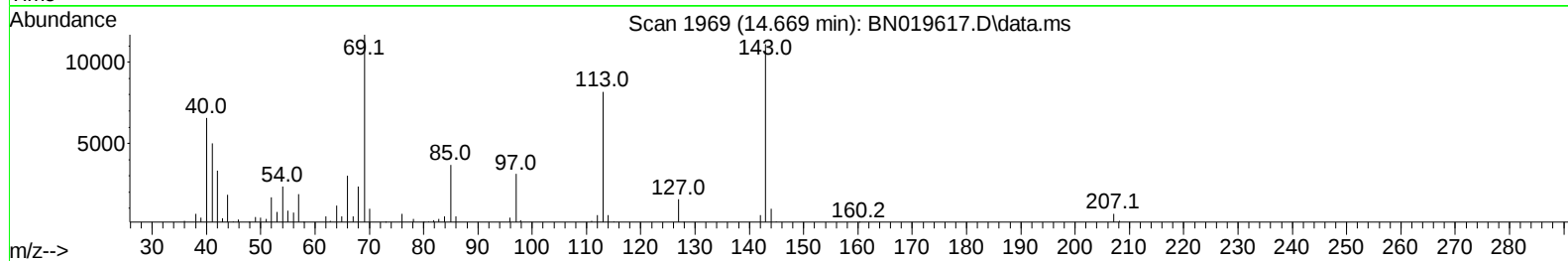
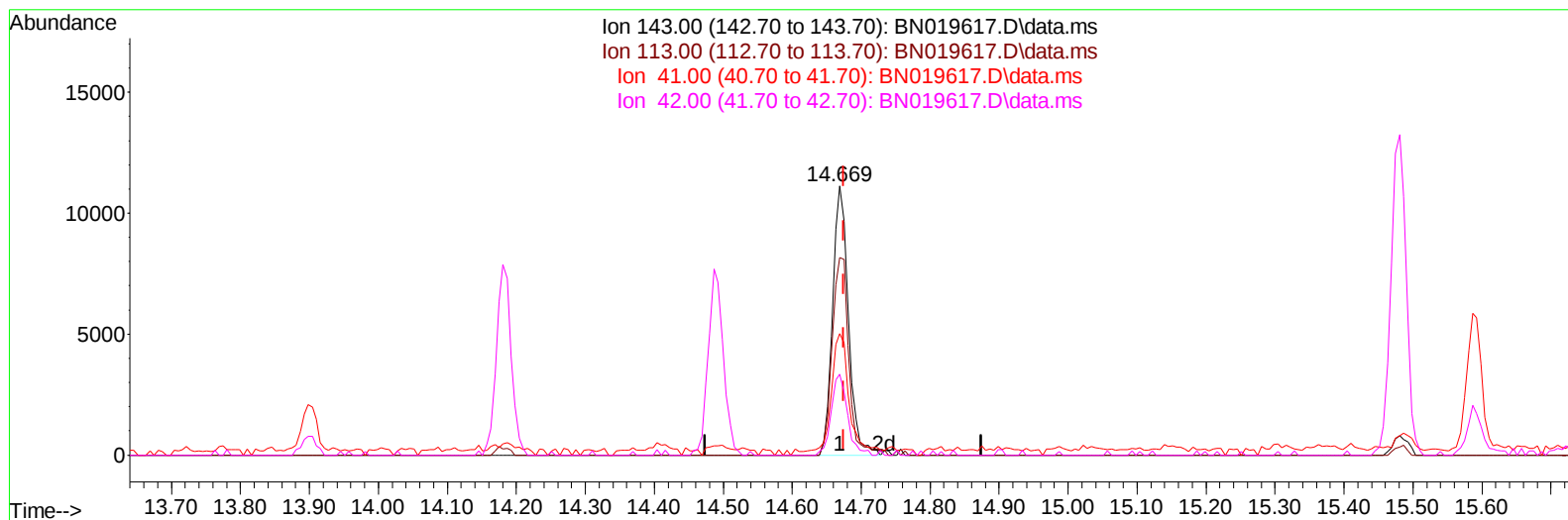
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Sample : N2621-01
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BFFH9

Manual IntegrationsAPPROVED

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TIC: BN019617.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (-0.006) 3.73 ng/u1 m

response 17599

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.20	73.55
41.00	39.70	45.09
42.00	27.00	30.04

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 ALS Vial : 7 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	155758	20.000	ng/ul	0.00
20) Naphthalene-d8	10.663	136	622283	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.492	164	353429	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.228	188	671264	20.000	ng/ul	0.00
79) Chrysene-d12	21.416	240	638657	20.000	ng/ul	0.00
88) Perylene-d12	23.768	264	659773	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	19497m	4.724	ng/uL	0.00
4) Pyridine-d5	3.752	84	53800	5.148	ng/ul	0.00
7) Phenol-d5	7.034	99	63908	4.735	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.199	67	237479	30.332	ng/ul	0.00
11) 2-Chlorophenol-d4	7.405	132	221690	20.706	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	126267	12.211	ng/ul	0.00
21) Nitrobenzene-d5	9.022	128	144201	31.782	ng/ul	0.00
24) 2-Nitrophenol-d4	9.746	143	126814	28.362	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	219000	24.468	ng/ul	0.00
31) 4-Chloroaniline-d4	10.793	131	285443	21.008	ng/ul	0.00
46) Dimethylphthalate-d6	13.898	166	804429	33.138	ng/ul	0.00
49) Acenaphthylene-d8	14.187	160	950925	28.972	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	17599m	3.728	ng/ul	0.00
60) Fluorene-d10	15.481	176	713047	33.274	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	77685	22.669	ng/ul	0.00
73) Anthracene-d10	17.328	188	1111146	35.149	ng/ul	0.00
81) Pyrene-d10	19.622	212	1281018	34.888	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.616	264	1216061	36.056	ng/ul	0.00

Target Compounds Qvalue

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

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