

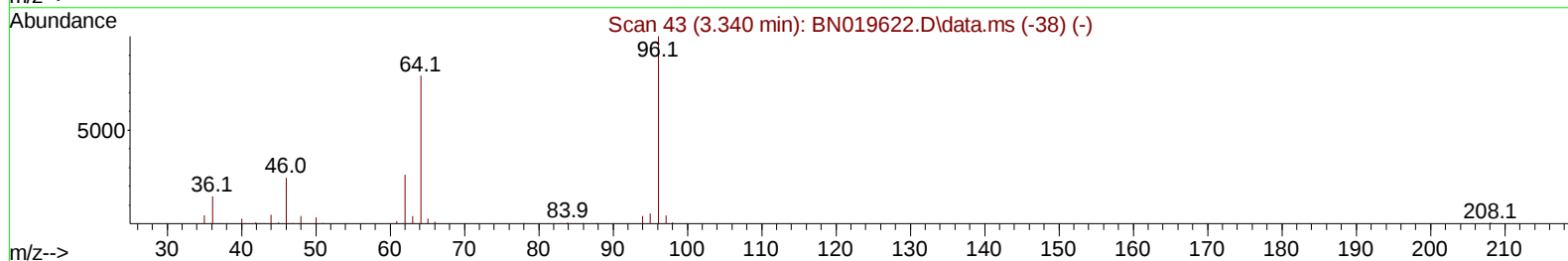
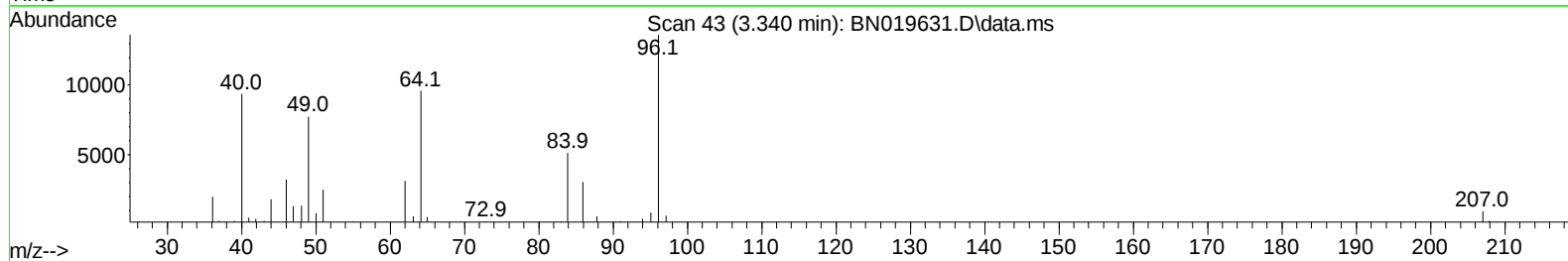
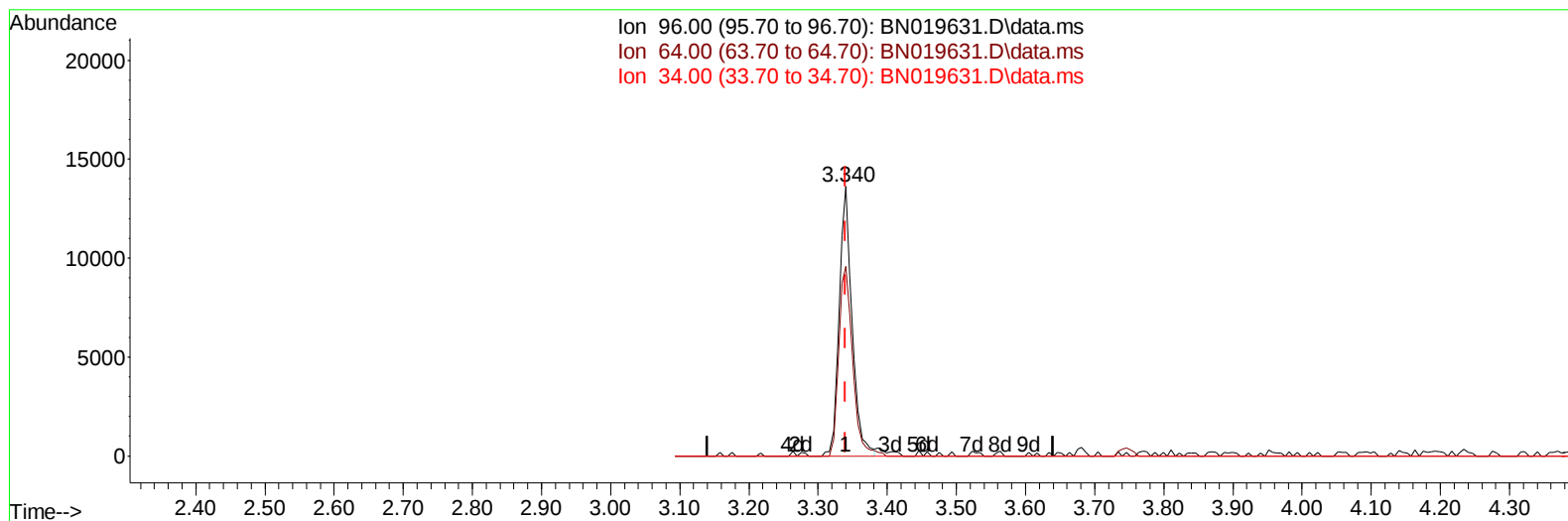
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN050222\
Data File : BN019631.D
Acq On : 02 May 2022 22:09
Operator : CG/JU
Sample : N2562-16
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
EW477

Manual IntegrationsAPPROVED

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

Quant Time: May 03 03:59:32 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



TIC: BN019631.D\data.ms

(3) 1,4-Dioxane-d8 (S)**3.340min (+ 0.000) 4.48 ng/uL****response 17753**

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	80.20	70.38
34.00	0.00	0.00
0.00	0.00	0.00

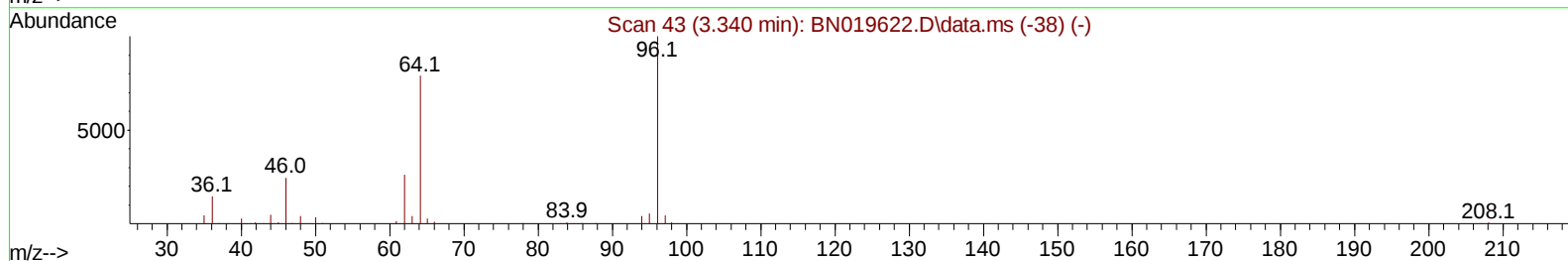
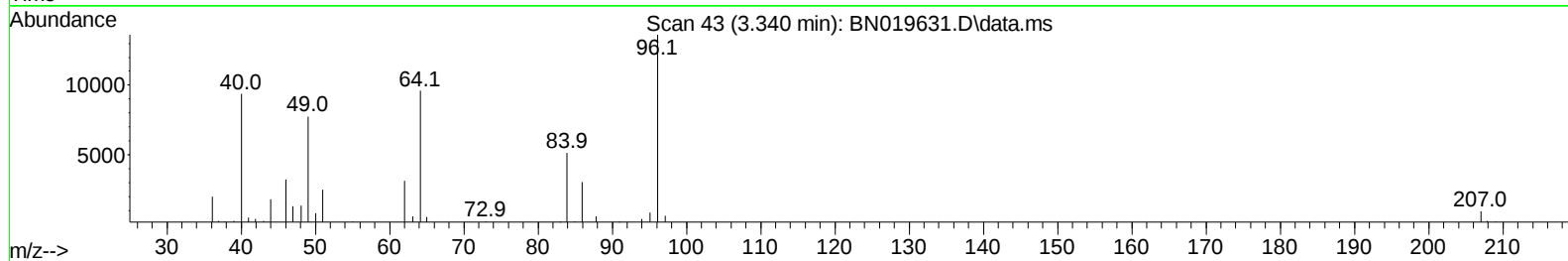
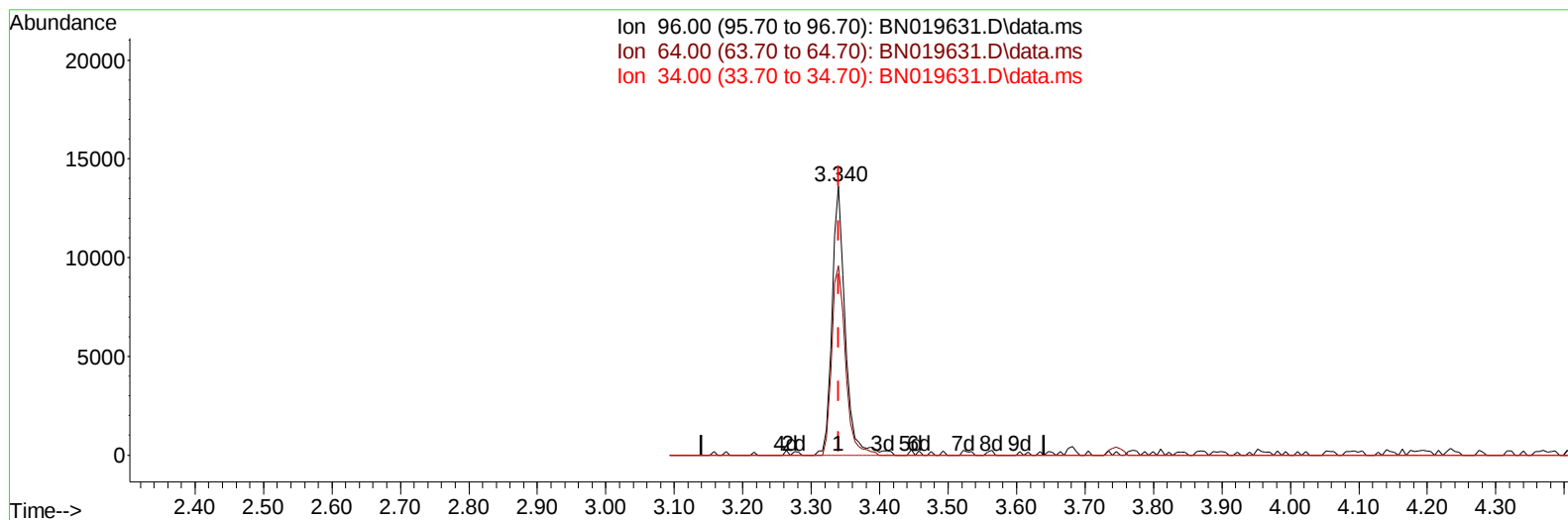
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TIC: BN019631.D\data.ms

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

3.340min (+ 0.000) 4.59 ng/uL m

response 18201

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	80.20	70.38
34.00	0.00	0.00
0.00	0.00	0.00

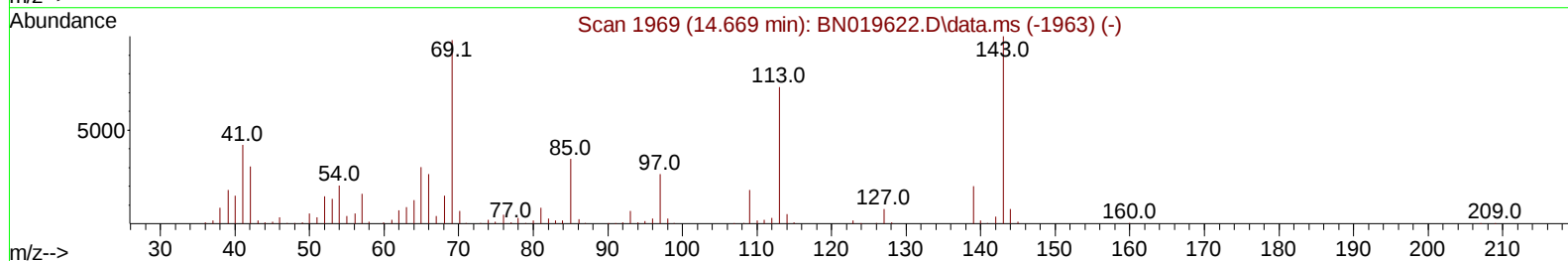
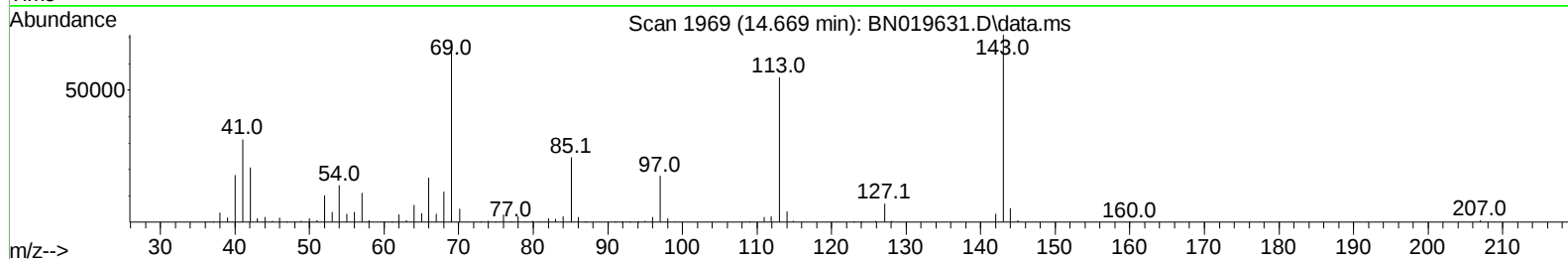
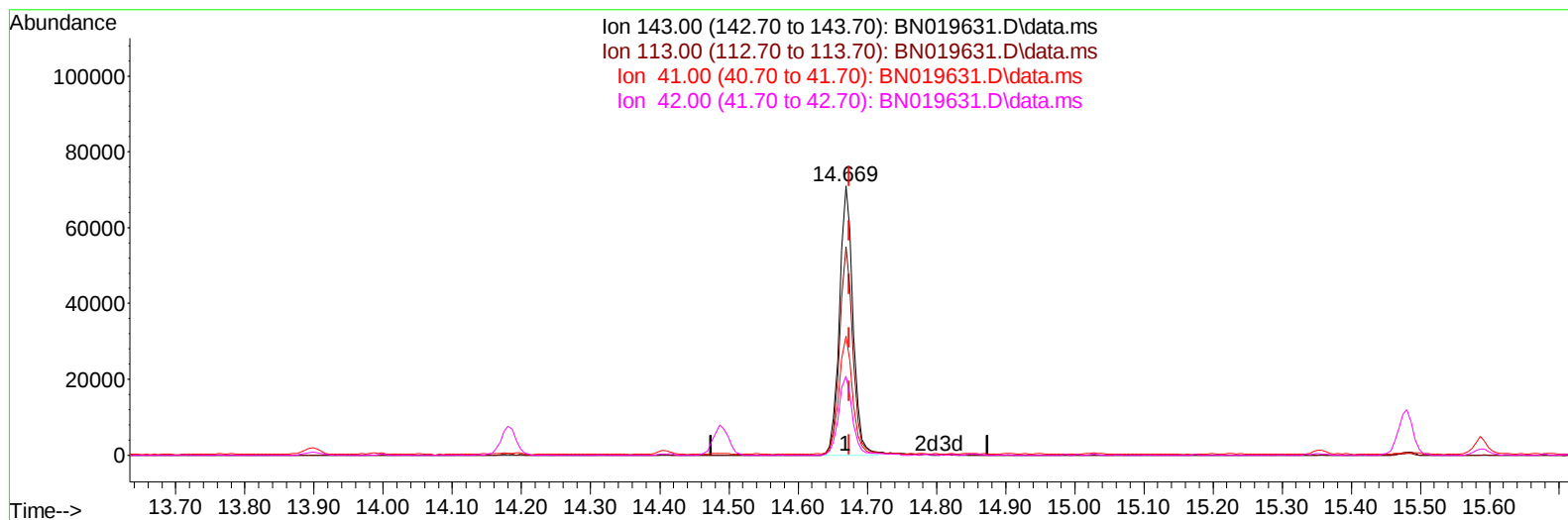
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Instrument :
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TIC: BN019631.D\data.ms

(54) 4-Nitrophenol-d4 (S)**14.669min (-0.006) 20.56 ng/u1****response 96879**

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.20	77.47
41.00	39.70	44.25
42.00	27.00	29.44

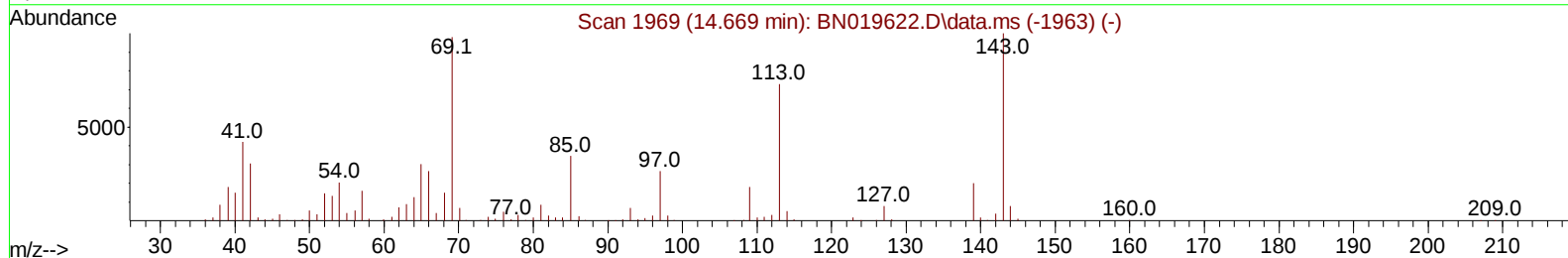
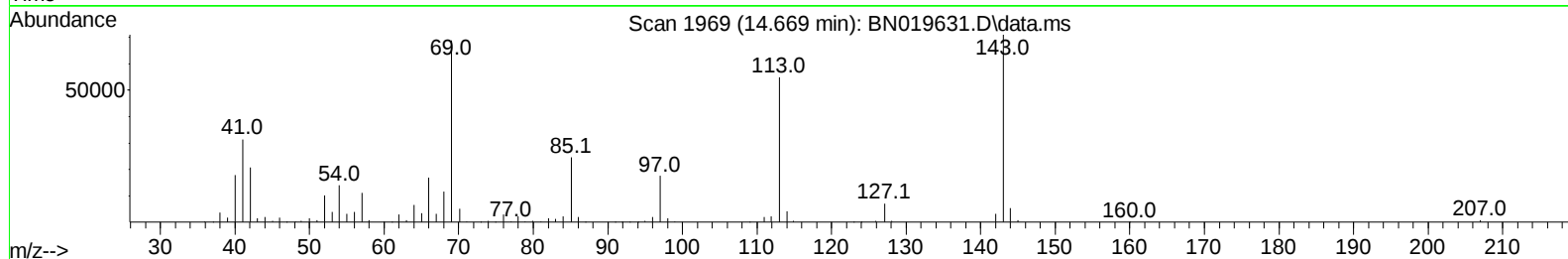
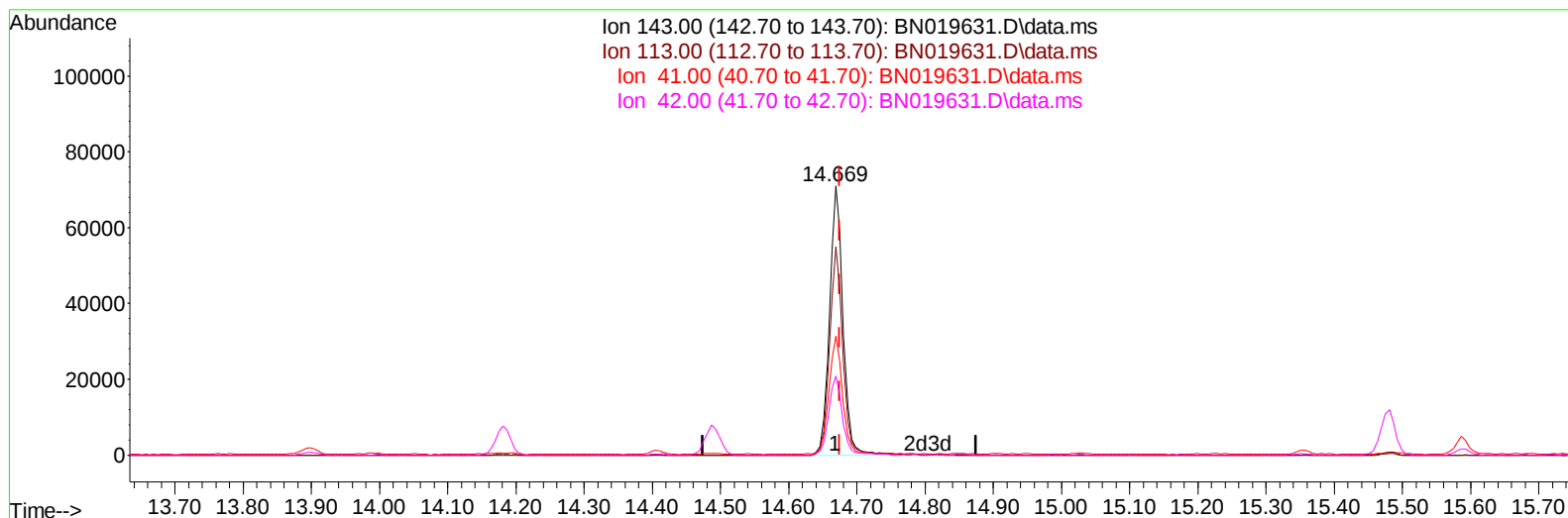
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Operator : CG/JU
Sample : N2562-16
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
EW477

Manual IntegrationsAPPROVED

Quant Time: May 03 03:59:32 2022
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Reviewed By :Jagrut Upadhyay 05/05/2022
Supervised By :Yogesh Patel 05/09/2022



TIC: BN019631.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (-0.006) 20.77 ng/ul m

response 97856

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.20	77.47
41.00	39.70	44.25
42.00	27.00	29.44

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	149686	20.000	ng/ul	0.00
20) Naphthalene-d8	10.663	136	606717	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.487	164	352760	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.228	188	692607	20.000	ng/ul	0.00
79) Chrysene-d12	21.410	240	671029	20.000	ng/ul	0.00
88) Perylene-d12	23.762	264	680190	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	18201m	4.589	ng/uL	0.00
4) Pyridine-d5	3.746	84	230256	22.924	ng/ul	0.00
7) Phenol-d5	7.028	99	292411	22.542	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.199	67	187800	24.960	ng/ul	0.00
11) 2-Chlorophenol-d4	7.399	132	248143	24.117	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	240930	24.246	ng/ul	0.00
21) Nitrobenzene-d5	9.016	128	119290	26.966	ng/ul	0.00
24) 2-Nitrophenol-d4	9.740	143	118818	27.255	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	213706	24.489	ng/ul	0.00
31) 4-Chloroaniline-d4	10.793	131	345188	26.057	ng/ul	0.00
46) Dimethylphthalate-d6	13.898	166	704605	29.081	ng/ul	0.00
49) Acenaphthylene-d8	14.181	160	853982	26.068	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	97856m	20.767	ng/ul	0.00
60) Fluorene-d10	15.481	176	620611	29.016	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.586	200	54746	15.483	ng/ul	0.00
73) Anthracene-d10	17.328	188	940436	28.833	ng/ul	0.00
81) Pyrene-d10	19.616	212	1088805	28.222	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.610	264	1011828	29.100	ng/ul	0.00

Target Compounds Qvalue

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

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