

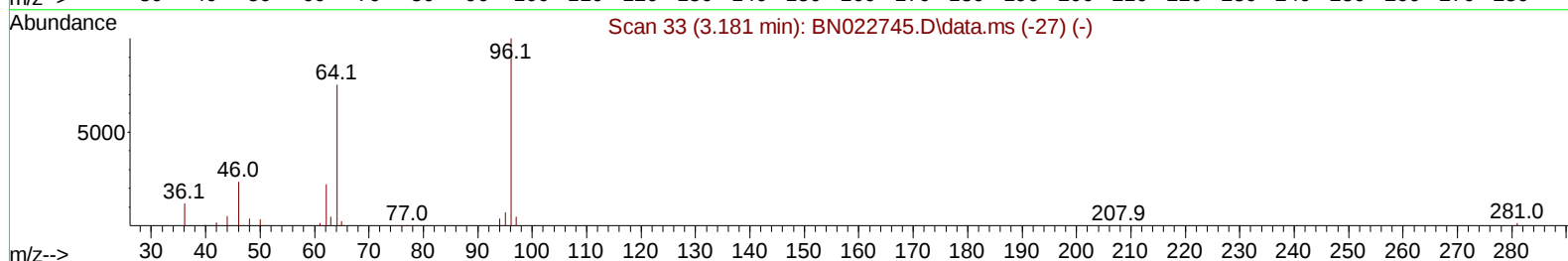
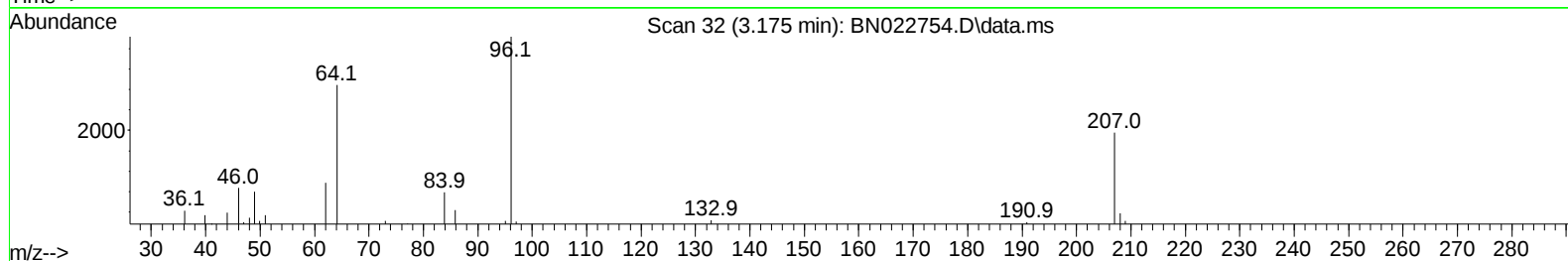
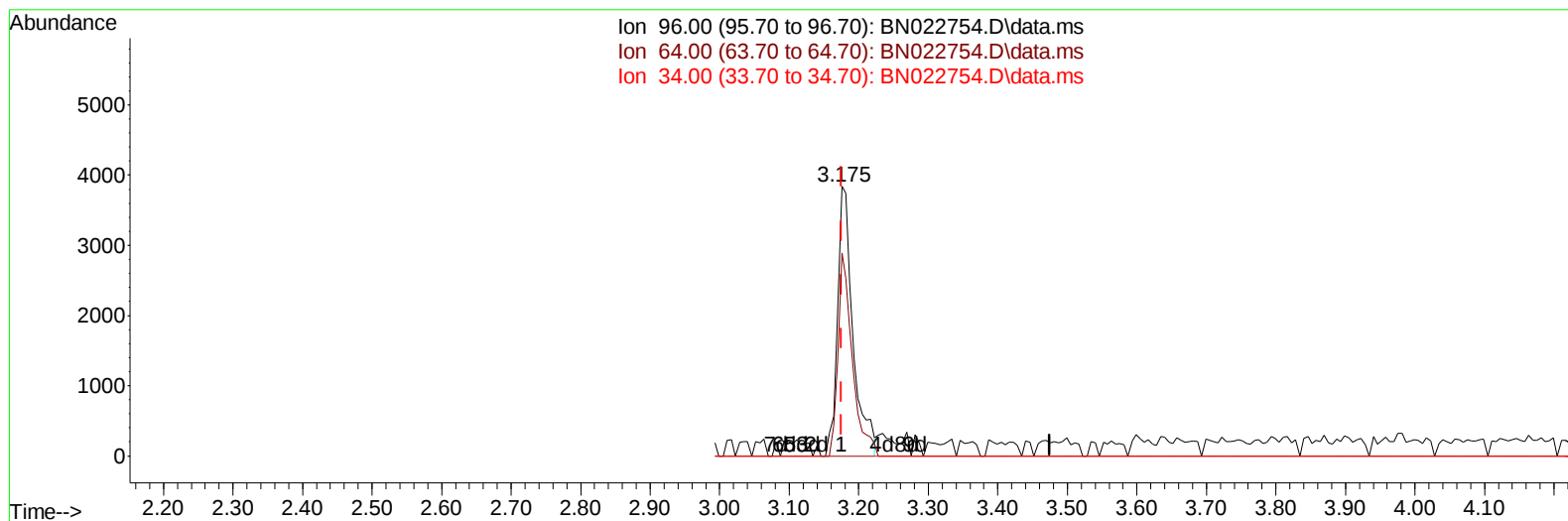
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111722\
 Data File : BN022754.D
 Acq On : 18 Nov 2022 13:23
 Operator : CG/JU
 Sample : N5689-02
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGCC5

Manual IntegrationsAPPROVED

Quant Time: Nov 18 21:48:11 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN111522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 18 05:17:17 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/21/2022
 Supervised By :mohammad ahmed 11/21/2022



TIC: BN022754.D\data.ms

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

3.175min (-0.000) 4.09 ng/uL

response 6023

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.50	75.33
34.00	0.00	0.00
0.00	0.00	0.00

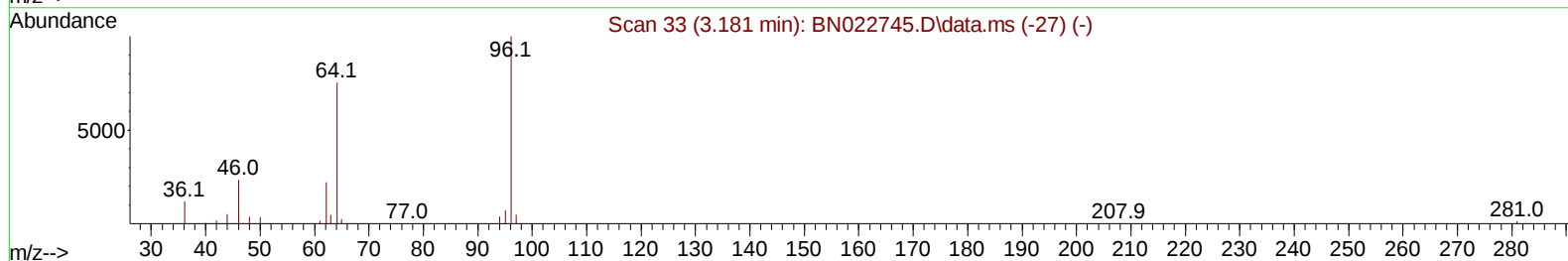
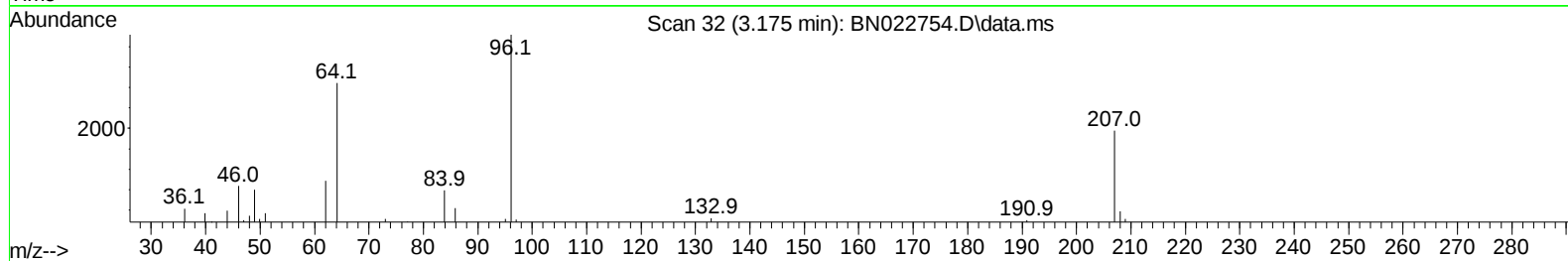
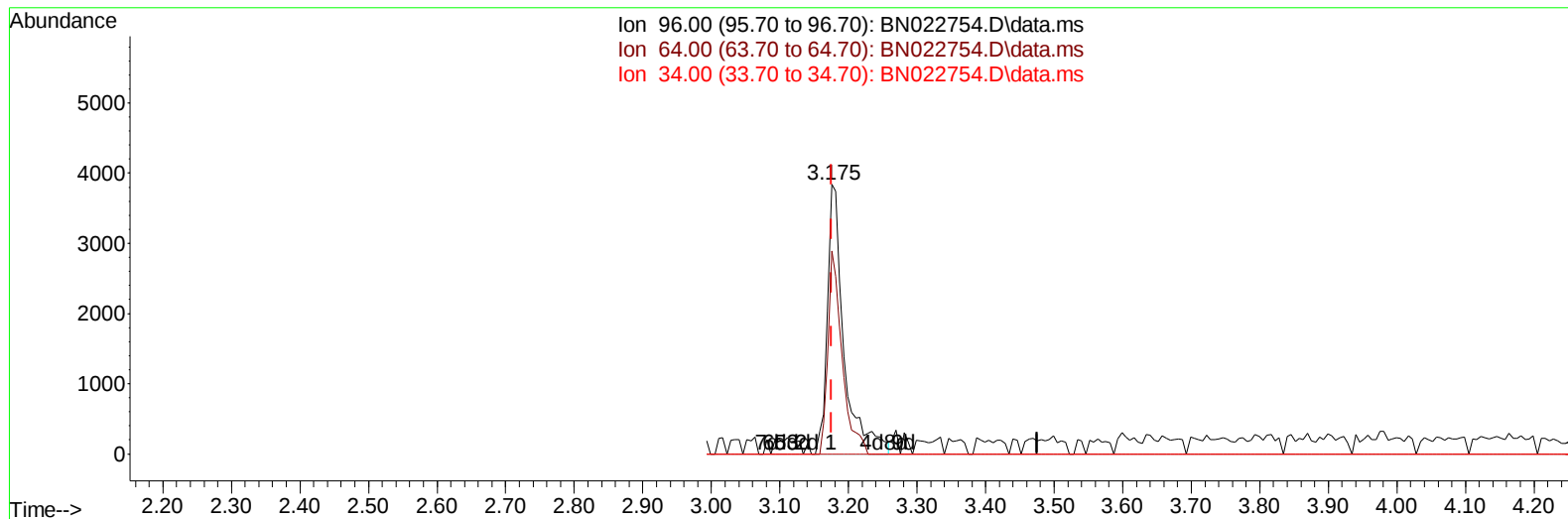
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TIC: BN022754.D\data.ms

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

3.175min (-0.000) 4.43 ng/uL m

response 6537

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.50	75.33
34.00	0.00	0.00
0.00	0.00	0.00

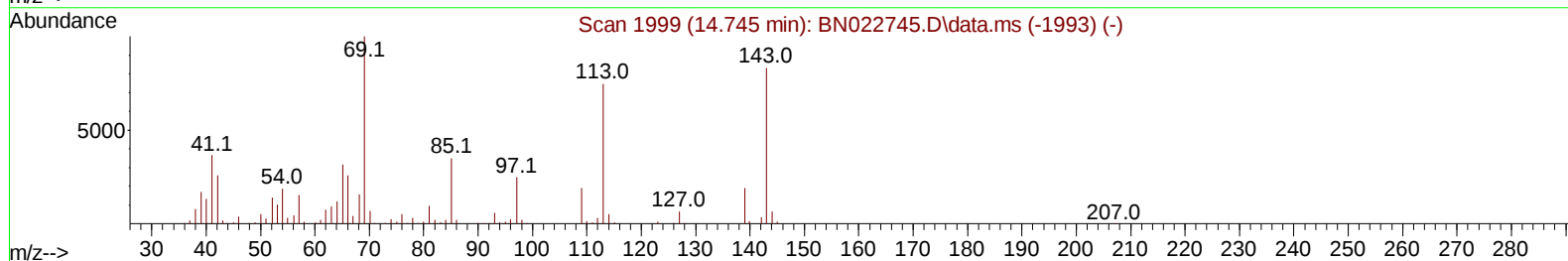
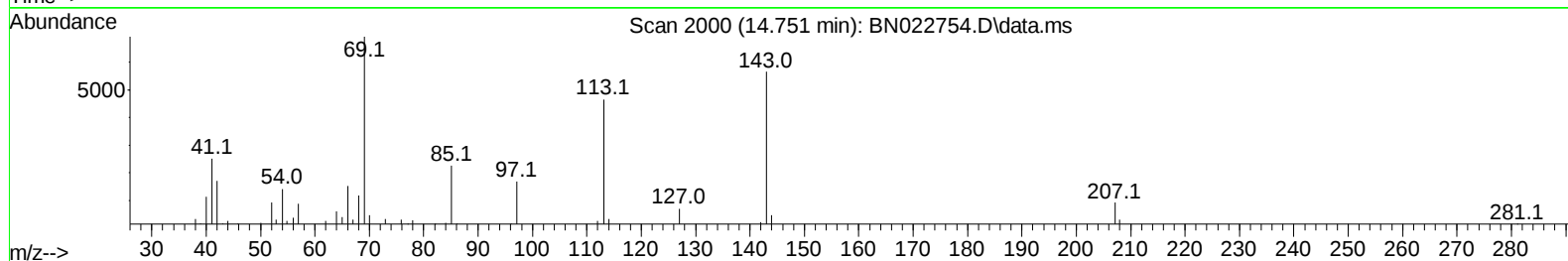
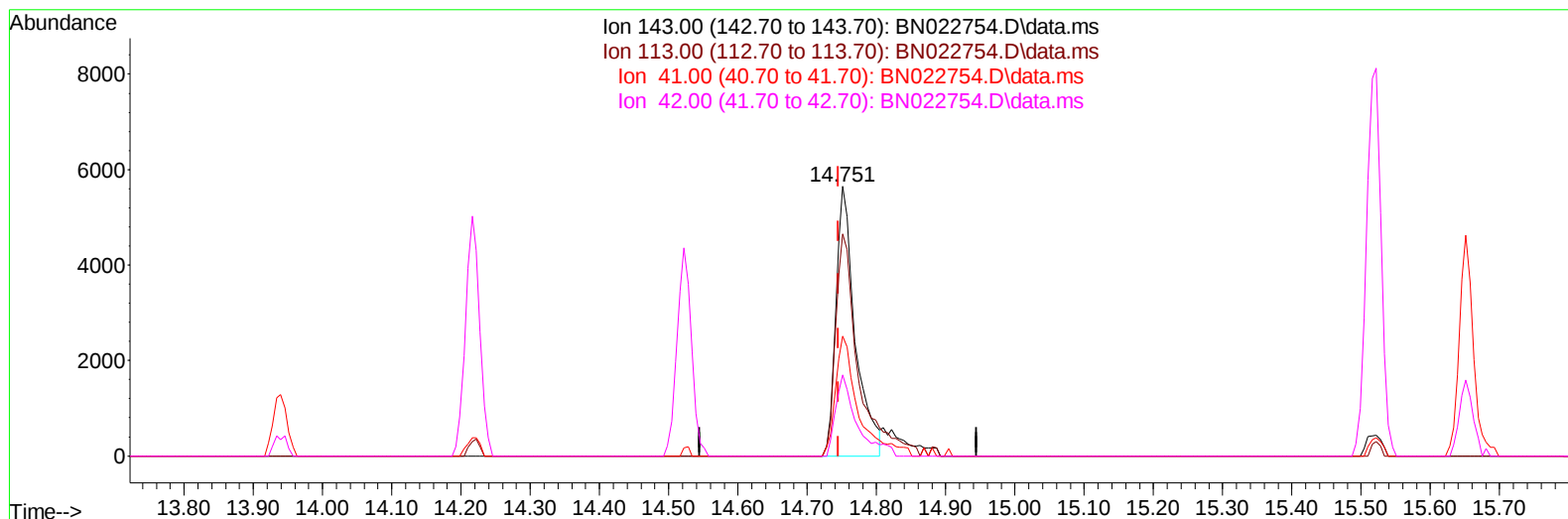
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Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
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TIC: BN022754.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (+ 0.006) 4.45 ng/u1

response 10881

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	95.00	82.47
41.00	46.30	44.47
42.00	31.30	30.13

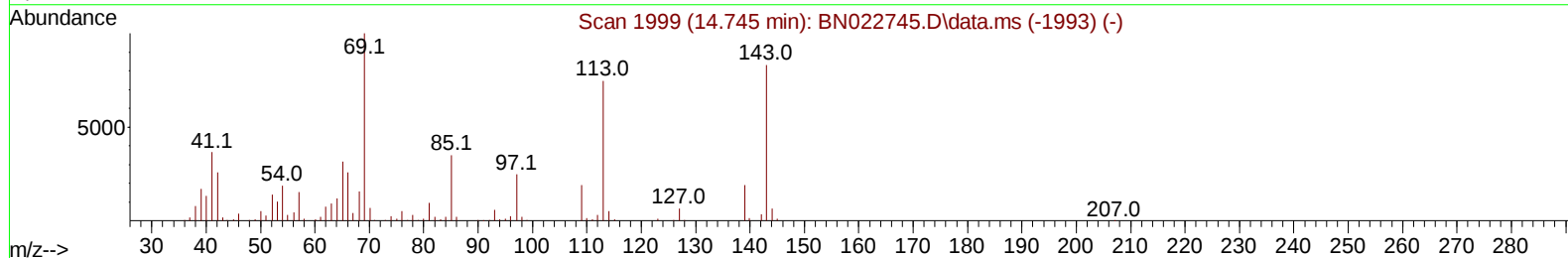
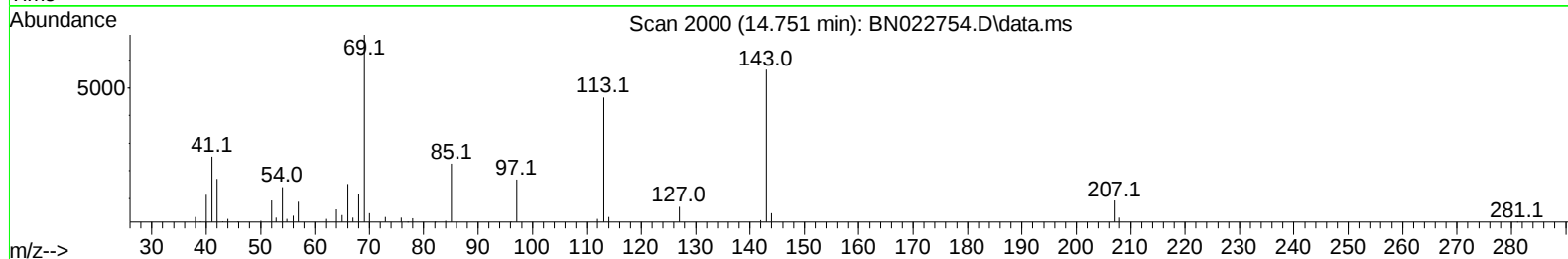
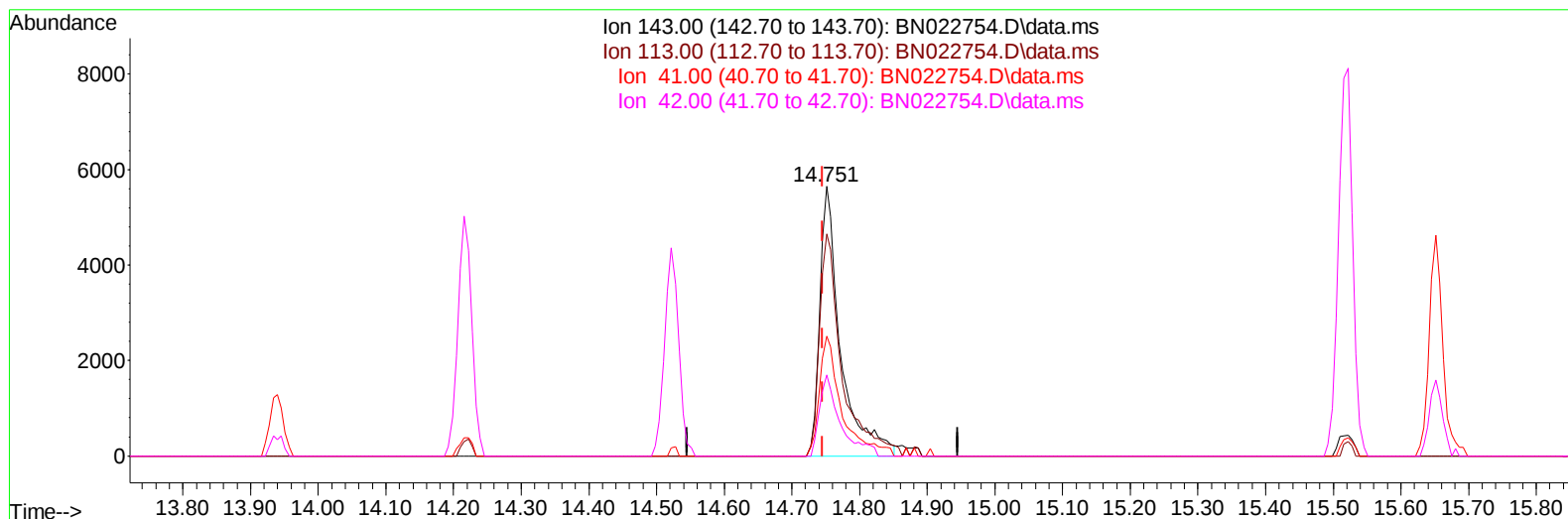
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Sample : N5689-02
Misc :
ALS Vial : 43 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
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Manual IntegrationsAPPROVED

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TIC: BN022754.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (+ 0.006) 4.90 ng/ul m

response 11978

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	95.00	82.47
41.00	46.30	44.47
42.00	31.30	30.13

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.852	152	50334	20.000	ng/ul	0.00
20) Naphthalene-d8	10.669	136	238780	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.522	164	168560	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.281	188	361629	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	374881	20.000	ng/ul	0.00
88) Perylene-d12	23.886	264	404322	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	6537m	4.434	ng/uL	0.00
4) Pyridine-d5	3.628	84	25897	6.211	ng/ul	0.01
7) Phenol-d5	7.022	99	28129	5.917	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	105902	37.154	ng/ul	0.00
11) 2-Chlorophenol-d4	7.381	132	87911	25.251	ng/ul	0.00
15) 4-Methylphenol-d8	8.575	113	59134	15.756	ng/ul	0.00
21) Nitrobenzene-d5	9.034	128	64707	35.010	ng/ul	0.00
24) 2-Nitrophenol-d4	9.757	143	64652	31.350	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.299	165	100008	28.445	ng/ul	0.00
31) 4-Chloroaniline-d4	10.822	131	140351	26.208	ng/ul	0.00
46) Dimethylphthalate-d6	13.940	166	450373	36.734	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	475938	35.231	ng/ul	0.00
54) 4-Nitrophenol-d4	14.751	143	11978m	4.901	ng/ul	0.00
60) Fluorene-d10	15.522	176	378103	37.602	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	60063	26.130	ng/ul	0.00
73) Anthracene-d10	17.381	188	628083	39.664	ng/ul	0.00
81) Pyrene-d10	19.680	212	725143	37.325	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.727	264	752651	37.031	ng/ul	0.00

Target Compounds Qvalue

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

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