

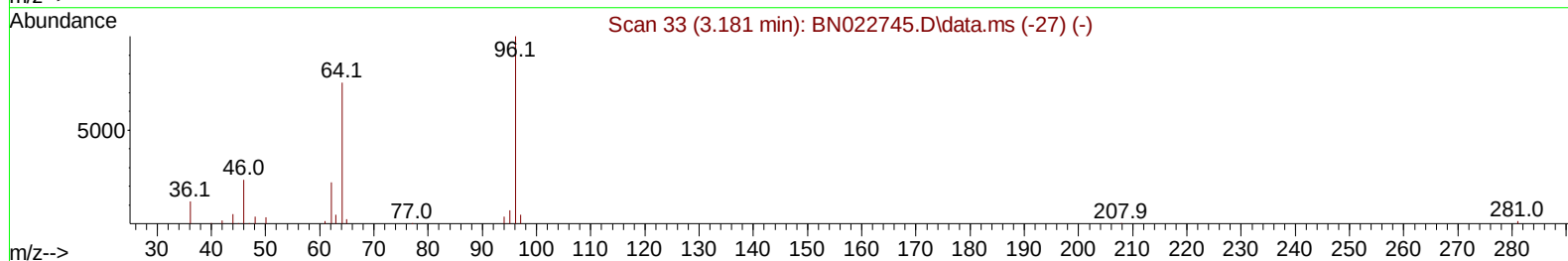
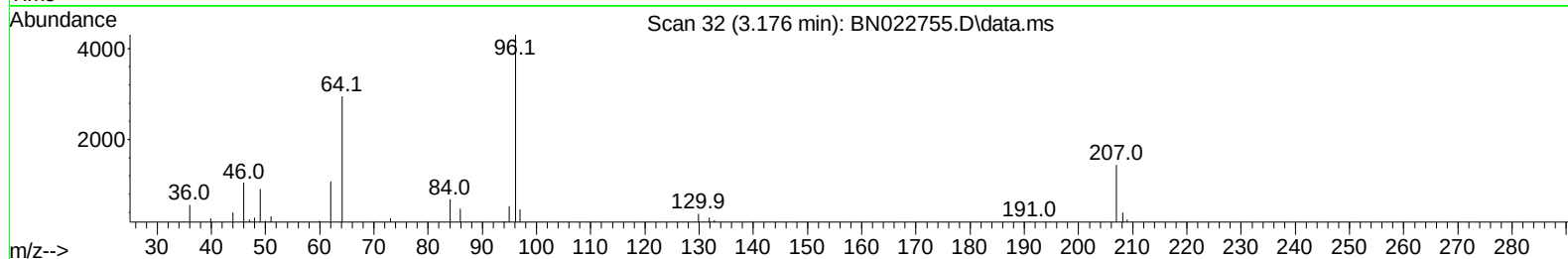
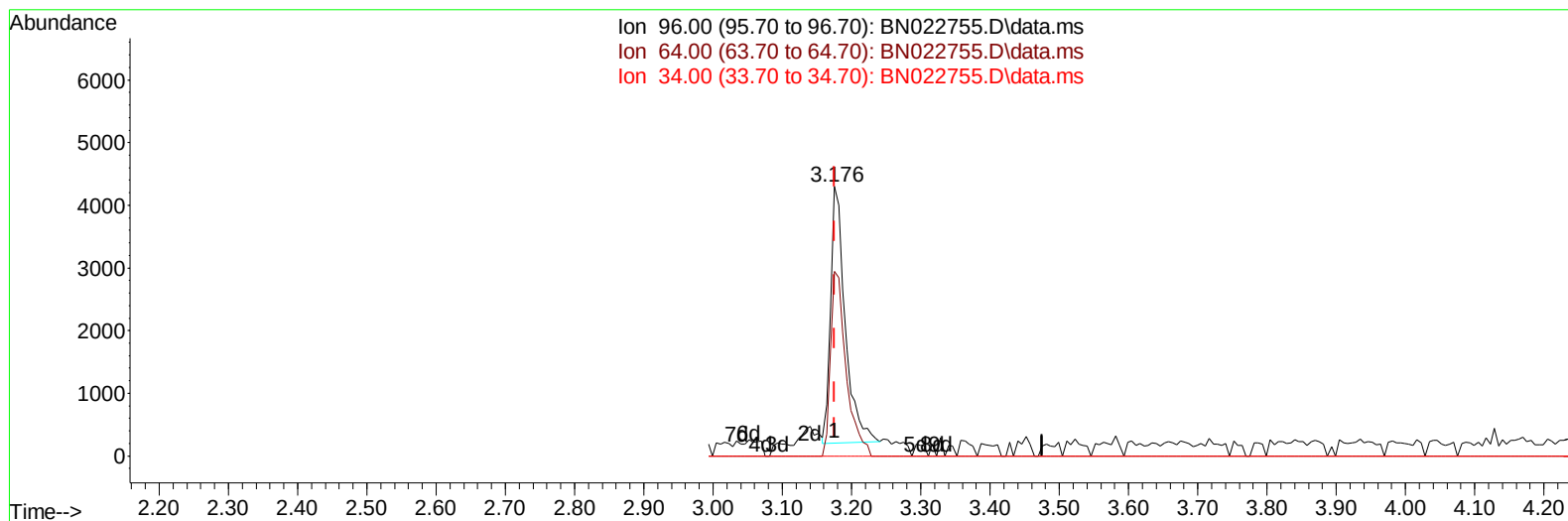
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111722\
 Data File : BN022755.D
 Acq On : 18 Nov 2022 14:00
 Operator : CG/JU
 Sample : N5689-10
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGCD3

Manual IntegrationsAPPROVED

Quant Time: Nov 18 21:48:18 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: BN022755.D\data.ms

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

3.176min (+ 0.000) 4.11 ng/uL

response 5984

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.50	68.60
34.00	0.00	0.00
0.00	0.00	0.00

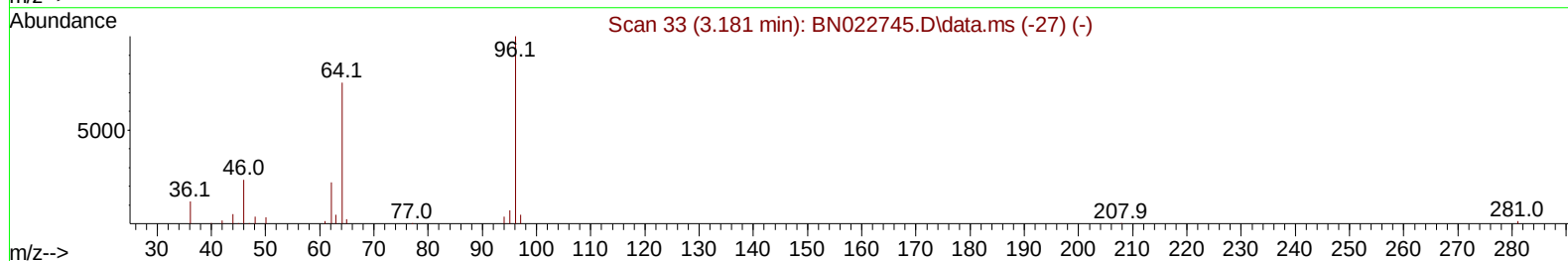
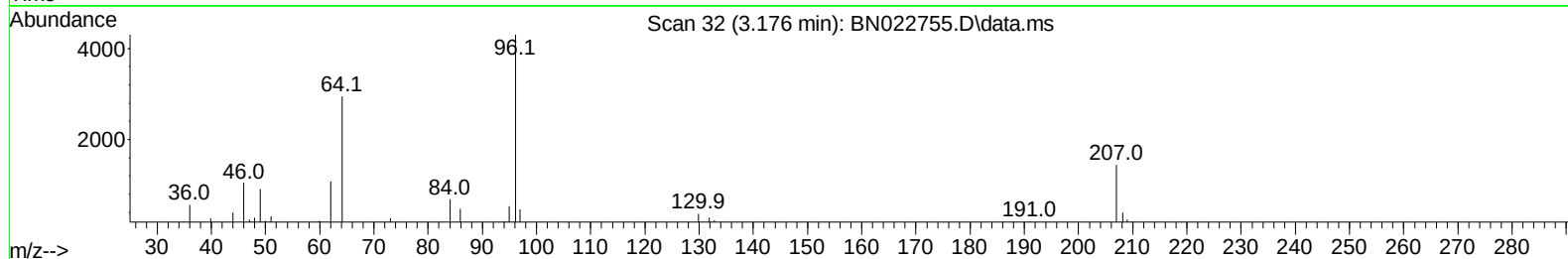
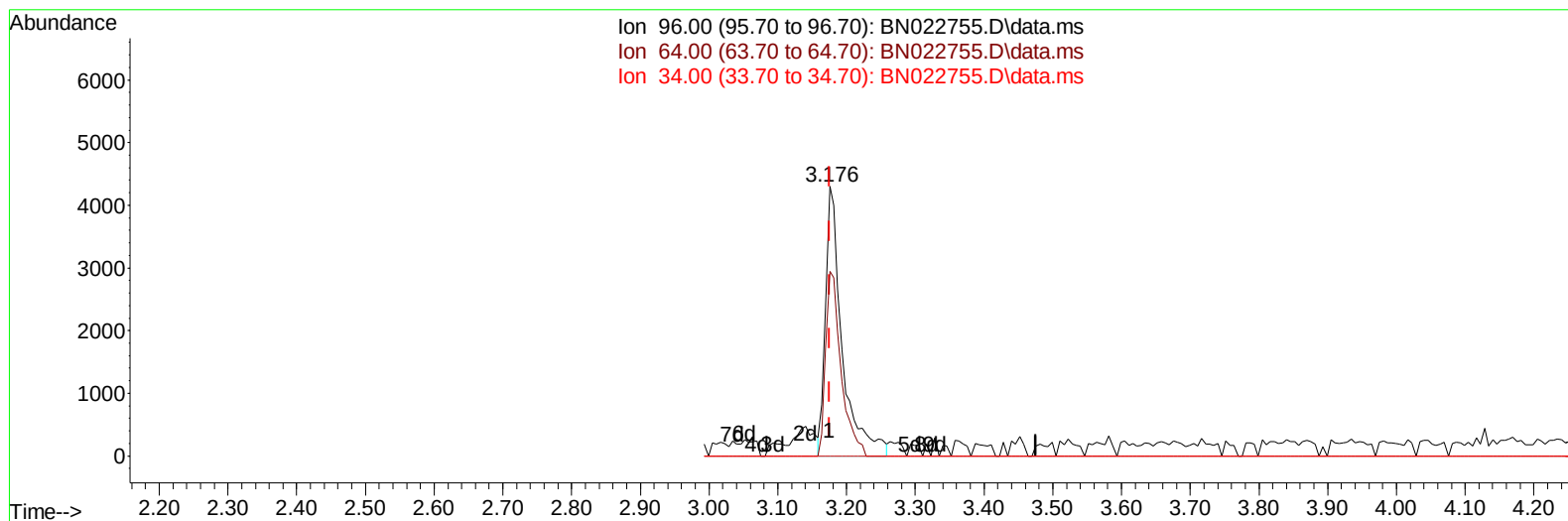
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(3) 1,4-Dioxane-d8 (S)

3.176min (+ 0.000) 5.02 ng/uL m

response 7313

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.50	68.60
34.00	0.00	0.00
0.00	0.00	0.00

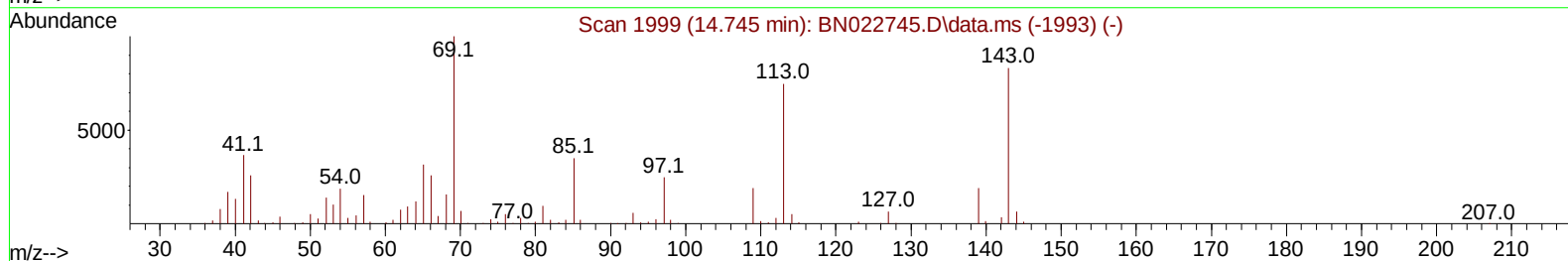
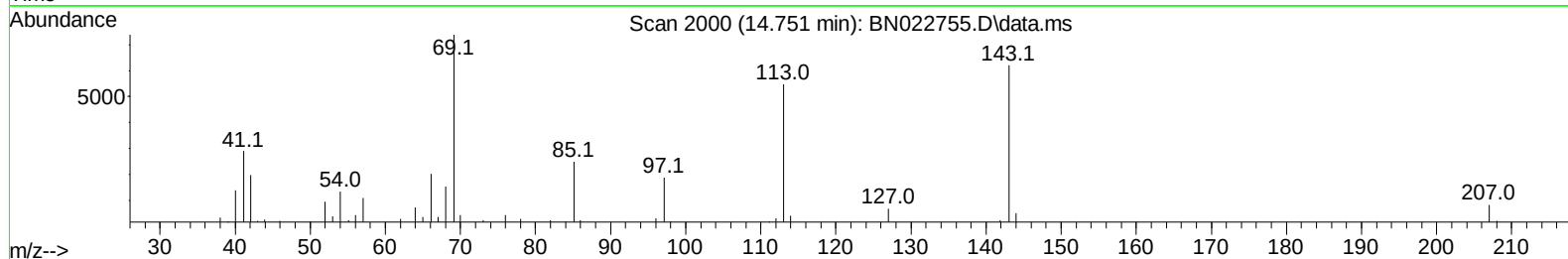
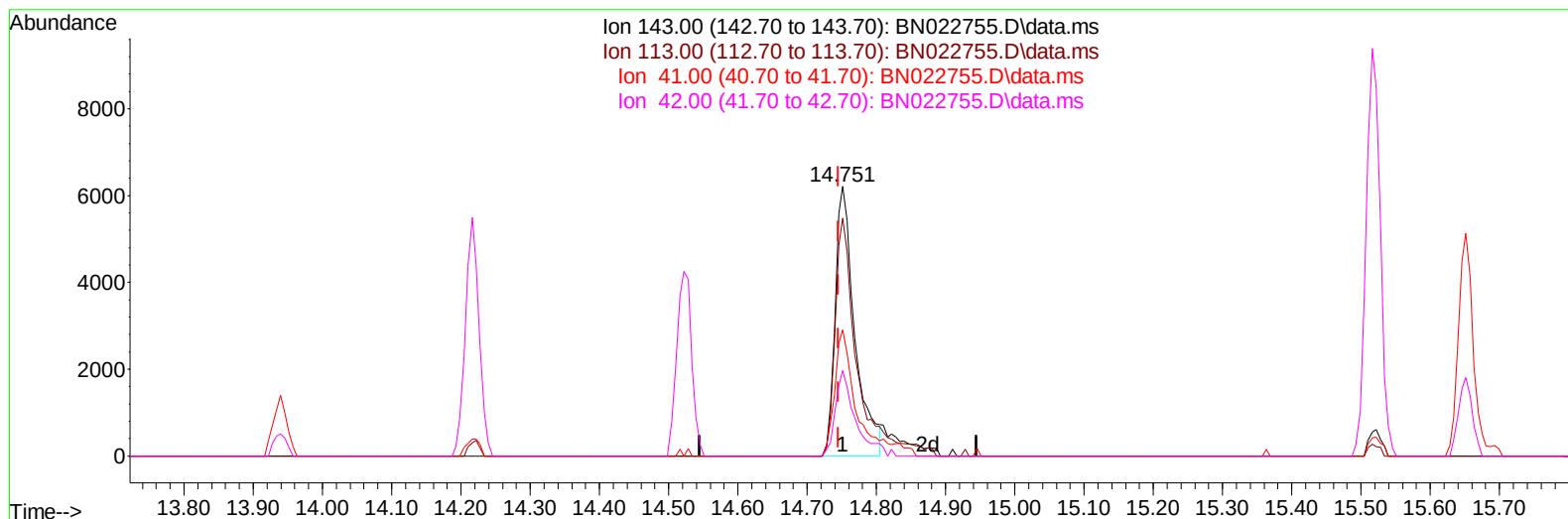
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Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
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TIC: BN022755.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (+ 0.006) 5.00 ng/ul

response 12231

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	95.00	88.34
41.00	46.30	46.74
42.00	31.30	31.68

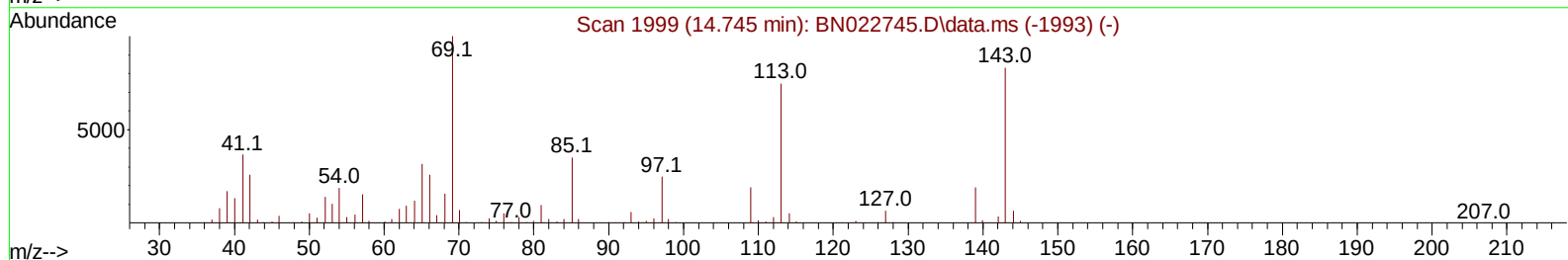
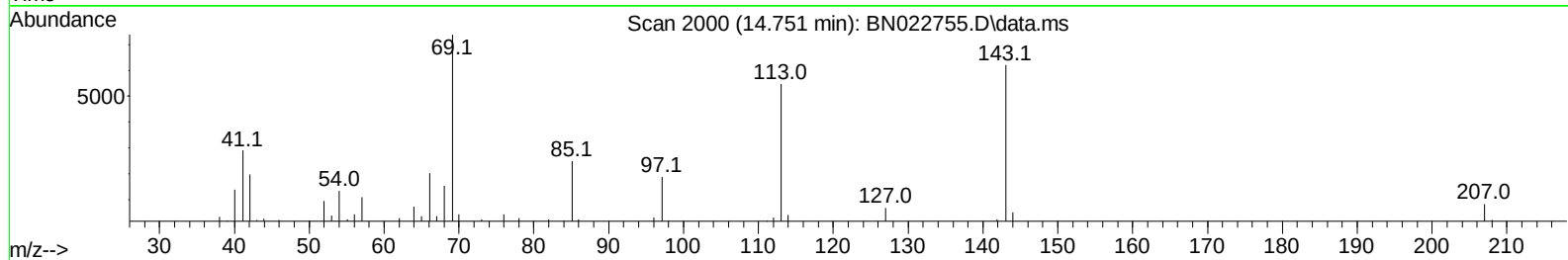
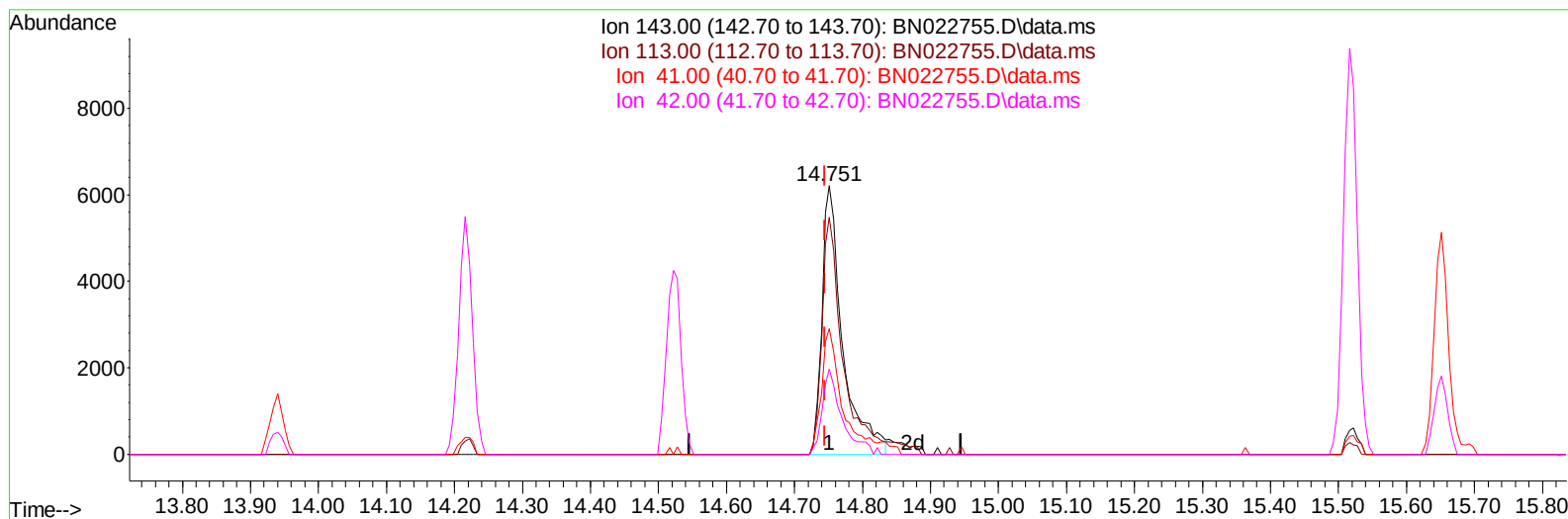
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Misc :
ALS Vial : 44 Sample Multiplier: 1

Instrument :
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TIC: BN022755.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (+ 0.006) 5.35 ng/ul m

response 13080

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	95.00	88.34
41.00	46.30	46.74
42.00	31.30	31.68

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.852	152	49732	20.000	ng/ul	0.00
20) Naphthalene-d8	10.669	136	241986	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.522	164	168733	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.275	188	362601	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	370788	20.000	ng/ul	0.00
88) Perylene-d12	23.886	264	402049	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.176	96	7313m	5.020	ng/uL	0.00
4) Pyridine-d5	3.634	84	18207	4.420	ng/ul	0.02
7) Phenol-d5	7.022	99	28875	6.147	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	108614	38.567	ng/ul	0.00
11) 2-Chlorophenol-d4	7.381	132	90434	26.290	ng/ul	0.00
15) 4-Methylphenol-d8	8.575	113	61137	16.487	ng/ul	0.00
21) Nitrobenzene-d5	9.034	128	67386	35.976	ng/ul	0.00
24) 2-Nitrophenol-d4	9.758	143	69759	33.378	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.293	165	108540	30.463	ng/ul	0.00
31) 4-Chloroaniline-d4	10.822	131	148312	27.328	ng/ul	0.00
46) Dimethylphthalate-d6	13.940	166	495599	40.381	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	514068	38.015	ng/ul	0.00
54) 4-Nitrophenol-d4	14.751	143	13080m	5.346	ng/ul	0.00
60) Fluorene-d10	15.522	176	410345	40.766	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	73119	31.725	ng/ul	0.00
73) Anthracene-d10	17.381	188	686965	43.267	ng/ul	0.00
81) Pyrene-d10	19.680	212	763608	39.739	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.727	264	825116	40.826	ng/ul	0.00

Target Compounds Qvalue

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

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