

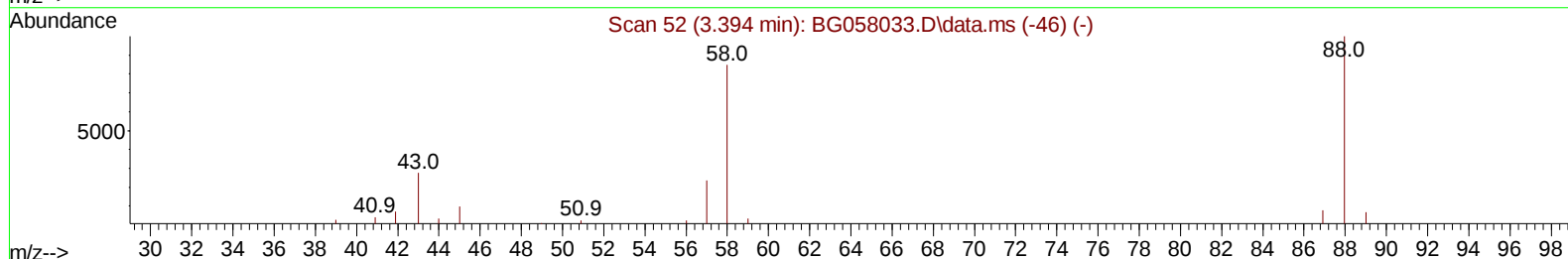
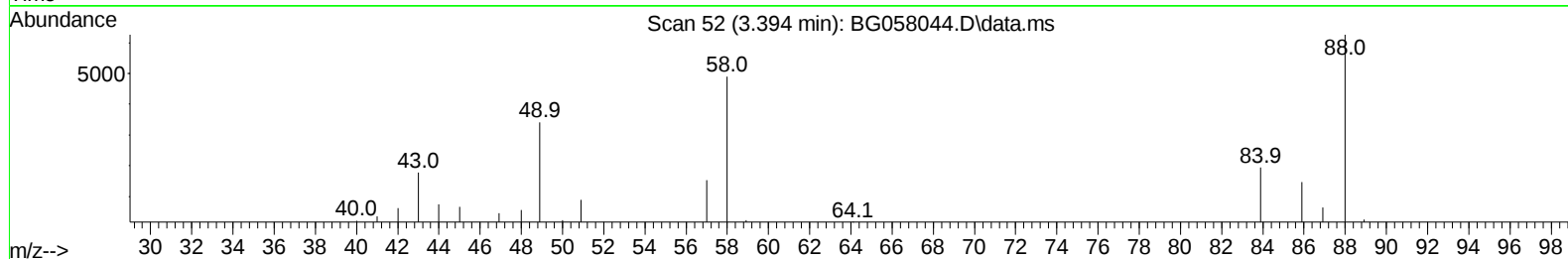
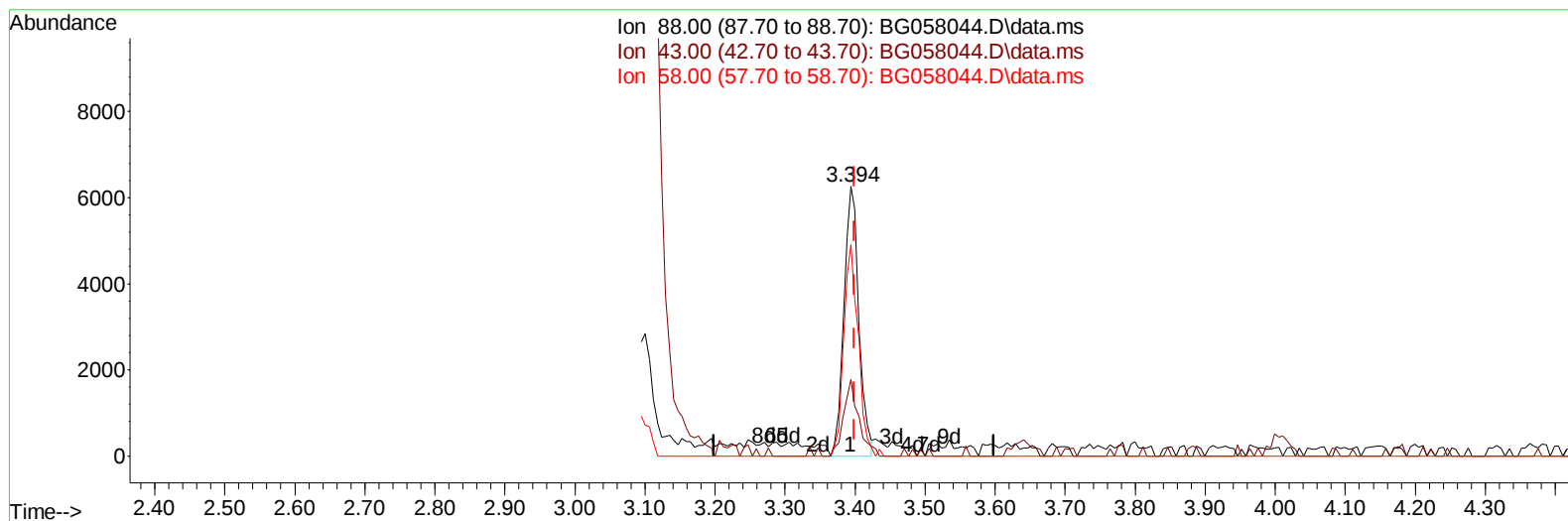
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070523\
 Data File : BG058044.D
 Acq On : 6 Jul 2023 10:44
 Operator : CG/JU
 Sample : 03258-07
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DCGW6

Manual IntegrationsAPPROVED

Quant Time: Jul 06 11:23:14 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 05 23:20:13 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 07/06/2023
 Supervised By :Sohil Jodhani 07/06/2023



TIC: BG058044.D\data.ms

(2) 1,4-Dioxane

3.394min (-0.005) 6.18 ng/uL

response 9409

Ion	Exp%	Act%
88.00	100.00	100.00
43.00	25.70	28.42
58.00	81.90	78.35
0.00	0.00	0.00

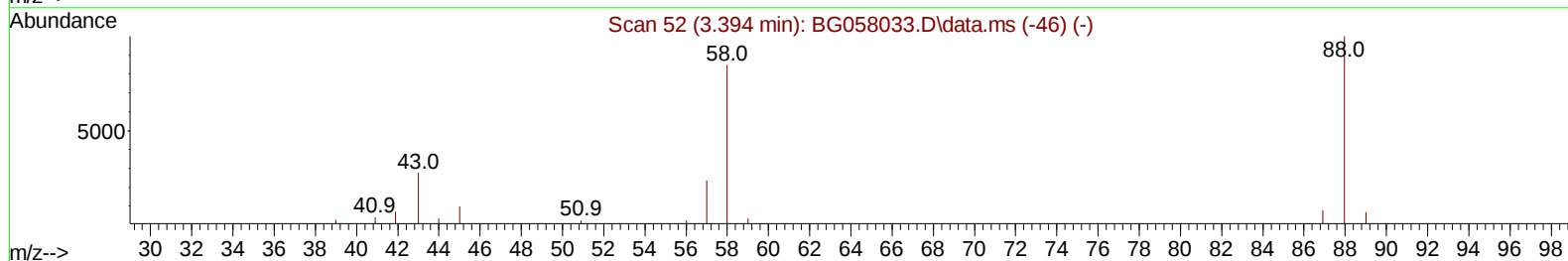
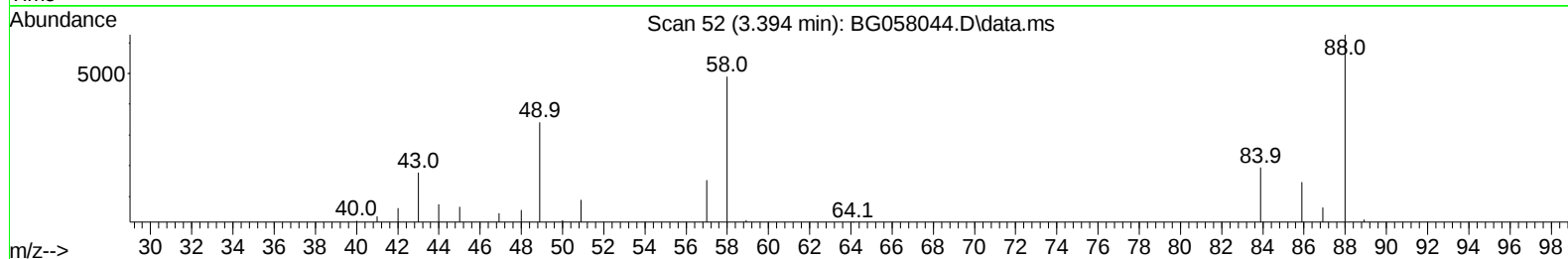
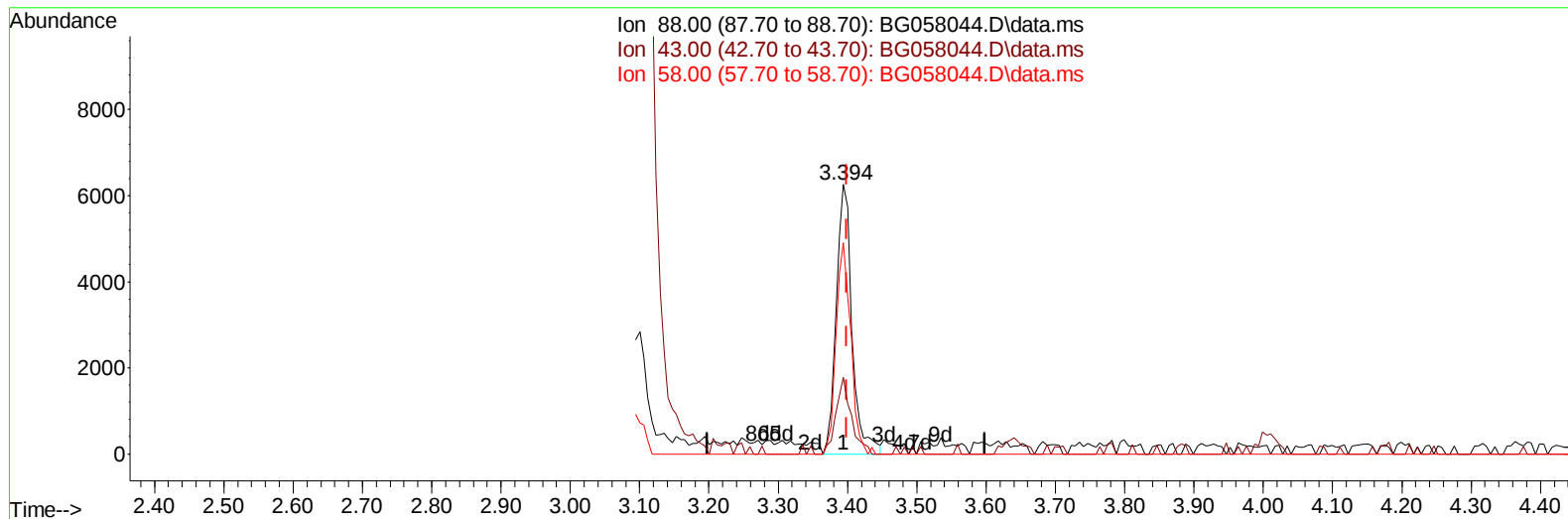
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TIC: BG058044.D\data.ms

(2) 1,4-Dioxane

3.394min (-0.005) 6.46 ng/uL m

response 9828

Ion	Exp%	Act%
88.00	100.00	100.00
43.00	25.70	28.42
58.00	81.90	78.35
0.00	0.00	0.00

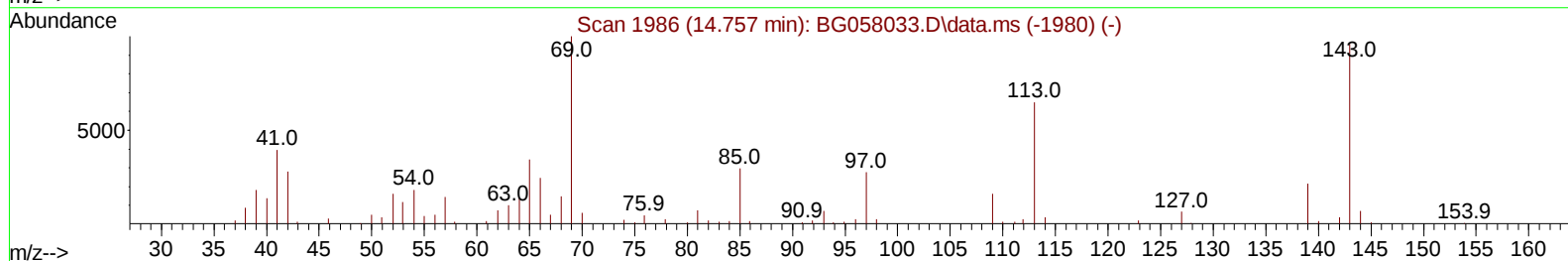
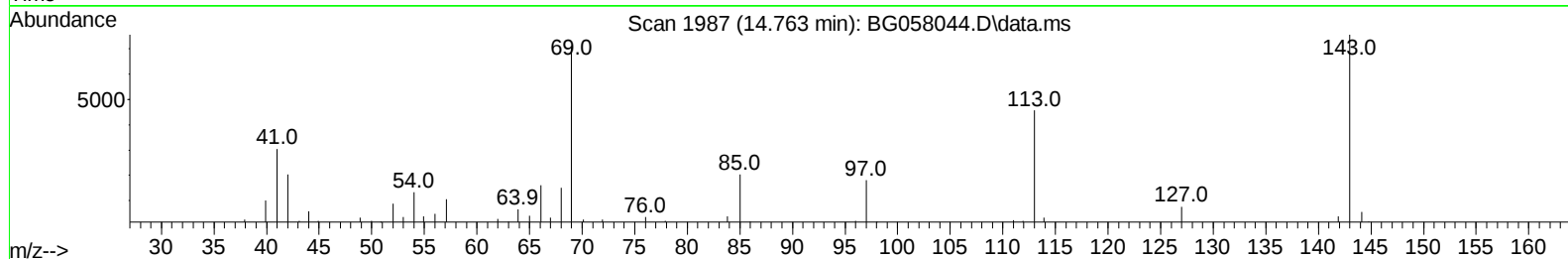
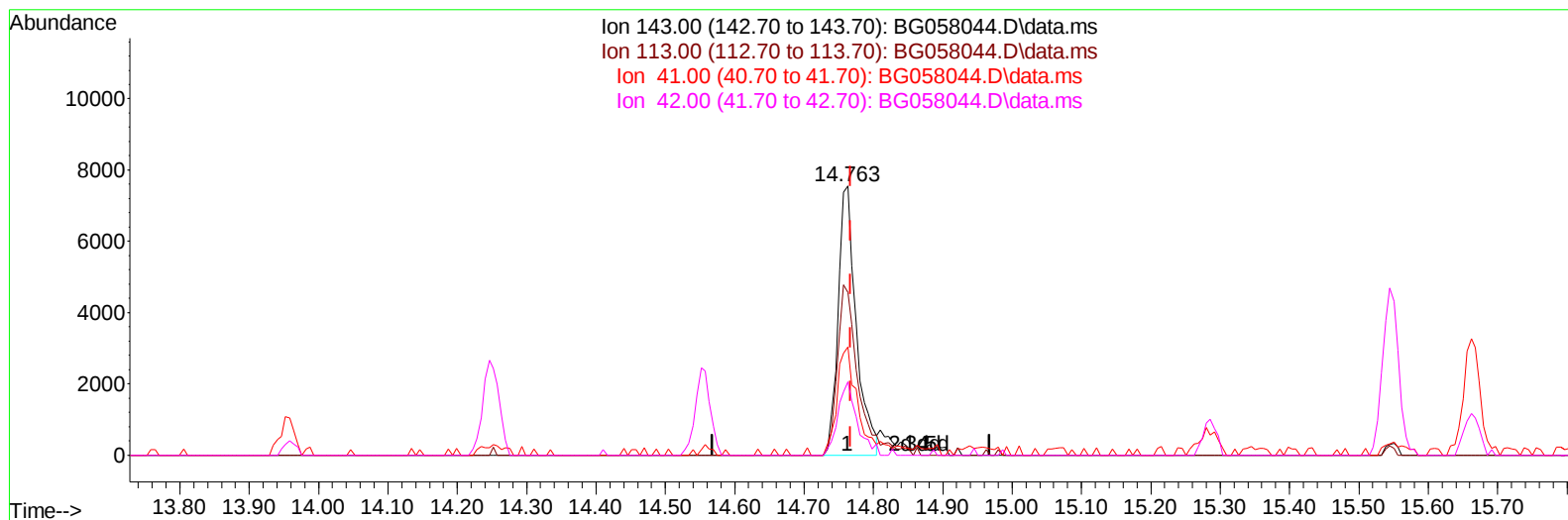
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ALS Vial : 28 Sample Multiplier: 1

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TIC: BG058044.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.005) 7.21 ng/u1

response 13695

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	66.40	60.50
41.00	40.50	40.28
42.00	27.50	27.15

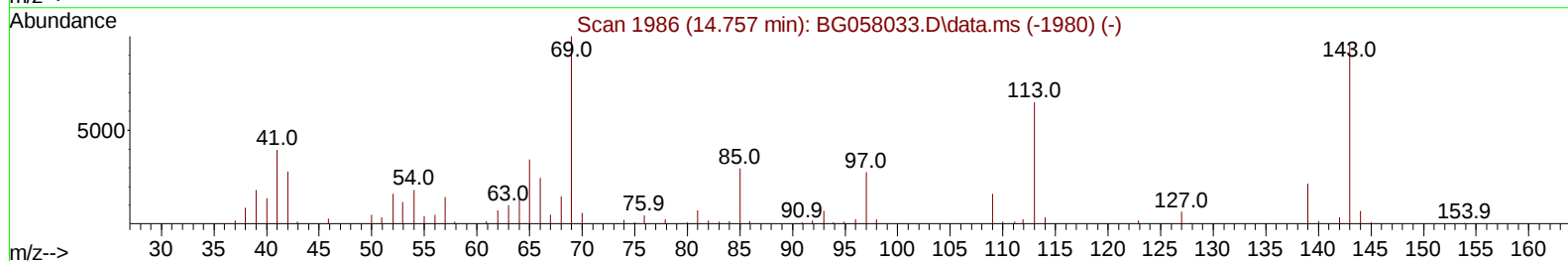
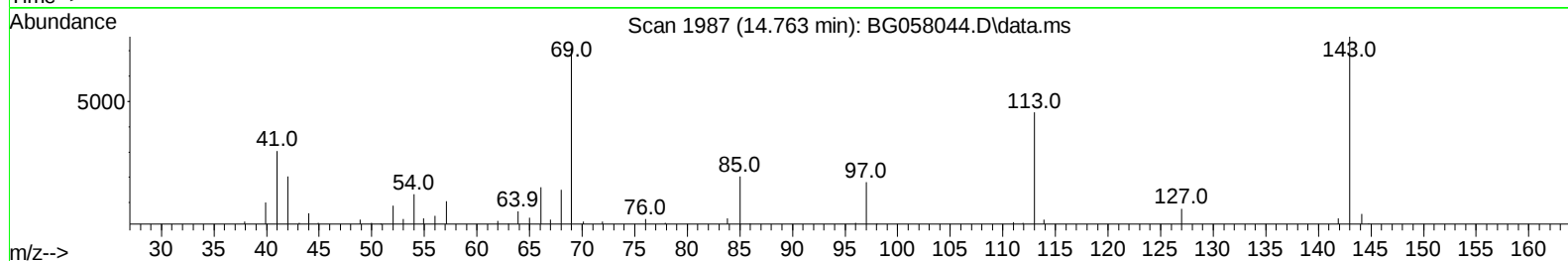
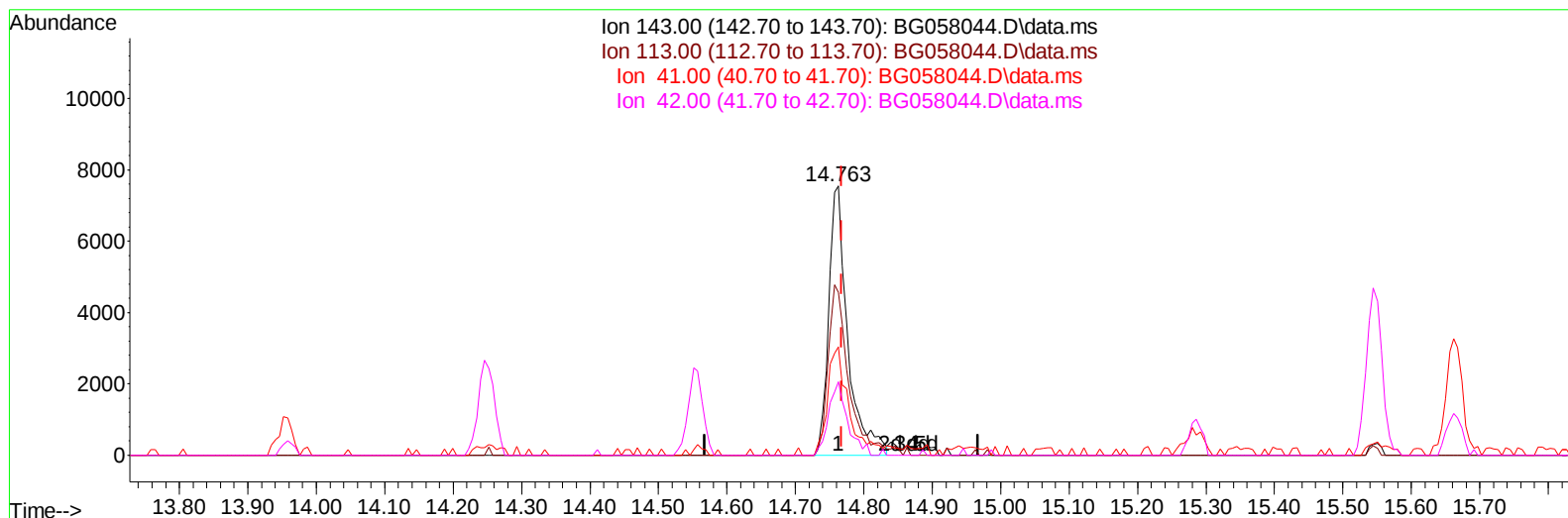
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ALS Vial : 28 Sample Multiplier: 1

Instrument :
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TIC: BG058044.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.005) 7.65 ng/ul m

response 14522

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	66.40	60.50
41.00	40.50	40.28
42.00	27.50	27.15

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/06/2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.918	152	43135	20.000	ng/ul	0.00
20) Naphthalene-d8	10.721	136	188225	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.557	164	136168	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.301	188	356890	20.000	ng/ul	0.00
79) Chrysene-d12	21.555	240	333545	20.000	ng/ul	0.00
88) Perylene-d12	24.716	264	415172	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.359	96	5822	4.726	ng/uL	0.00
4) Pyridine-d5	3.770	84	38969	10.420	ng/ul	0.00
7) Phenol-d5	7.084	99	31636	7.182	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.242	67	86740	32.674	ng/ul	0.00
11) 2-Chlorophenol-d4	7.448	132	78819	25.942	ng/ul	-0.01
15) 4-Methylphenol-d8	8.623	113	56274	16.930	ng/ul	-0.01
21) Nitrobenzene-d5	9.081	128	52860	34.456	ng/ul	0.00
24) 2-Nitrophenol-d4	9.804	143	54846	34.016	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.339	165	95627	30.917	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.856	131	116605	24.865	ng/ul	0.00
46) Dimethylphthalate-d6	13.958	166	382676	35.618	ng/ul	0.00
49) Acenaphthylene-d8	14.252	160	385833	32.910	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	14522m	7.651	ng/ul	0.00
60) Fluorene-d10	15.544	176	327820	35.928	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.662	200	64056	33.054	ng/ul	-0.01
73) Anthracene-d10	17.401	188	585441	35.351	ng/ul	0.00
81) Pyrene-d10	19.687	212	749860	35.696	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.499	264	800585	37.849	ng/ul	0.00
Target Compounds						
2) 1,4-Dioxane	3.394	88	9828m	6.458	ng/uL	Qvalue

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

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