

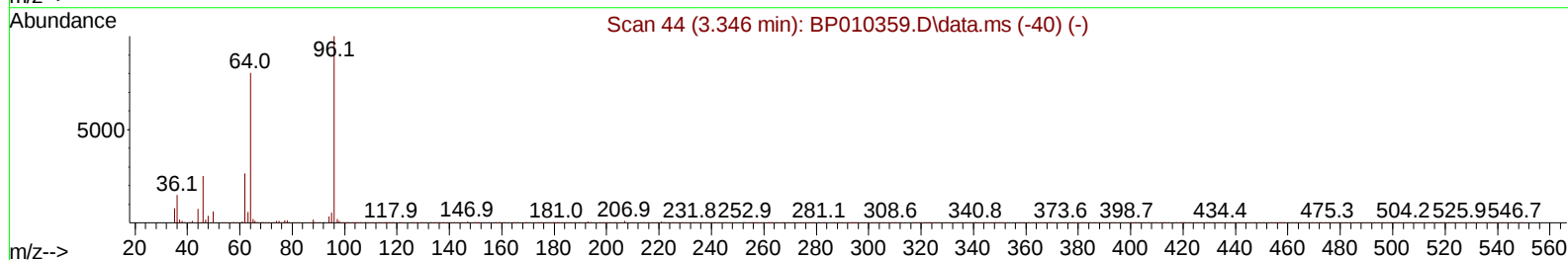
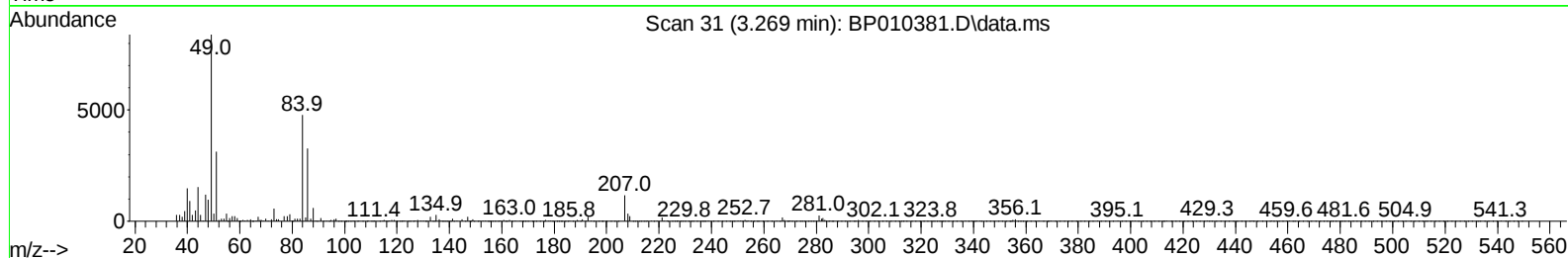
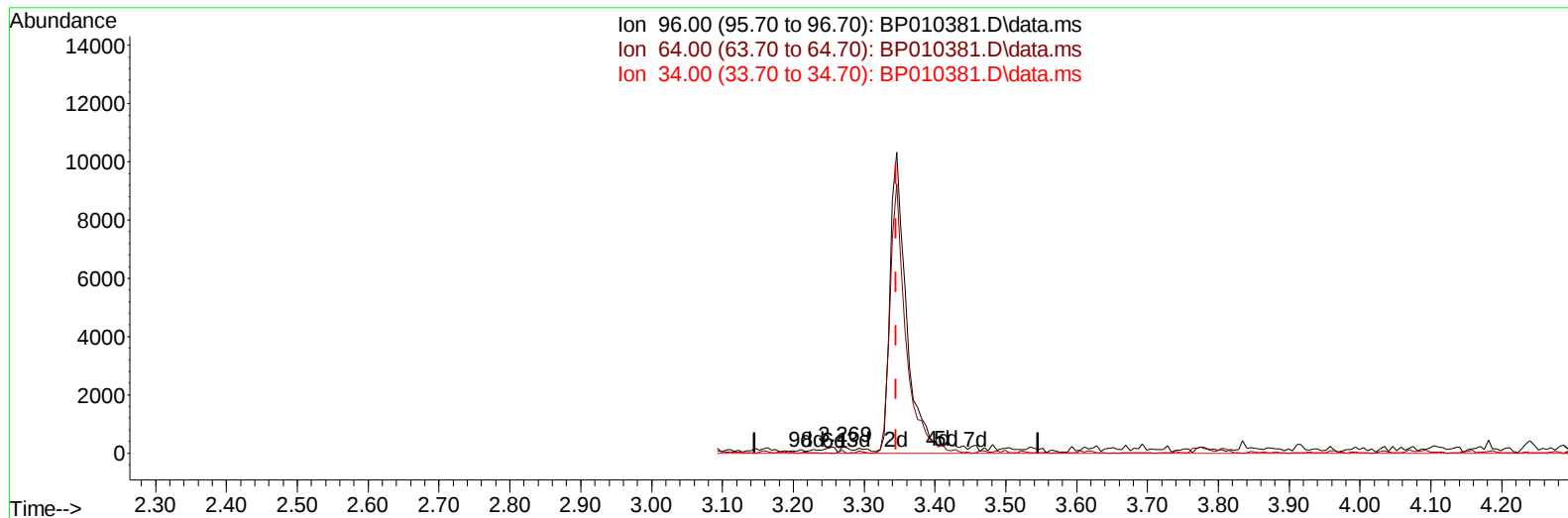
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051622\
 Data File : BP010381.D
 Acq On : 17 May 2022 05:22
 Operator : CG/JU
 Sample : N2831-07
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 YB2L6

Manual IntegrationsAPPROVED

Quant Time: May 17 07:07:59 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 13 15:04:31 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/17/2022
 Supervised By :mohammad ahmed 05/18/2022



TIC: BP010381.D\data.ms

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

3.269min (-0.077) 0.04 ng/uL

response 163

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	52.86
34.00	0.00	0.00
0.00	0.00	0.00

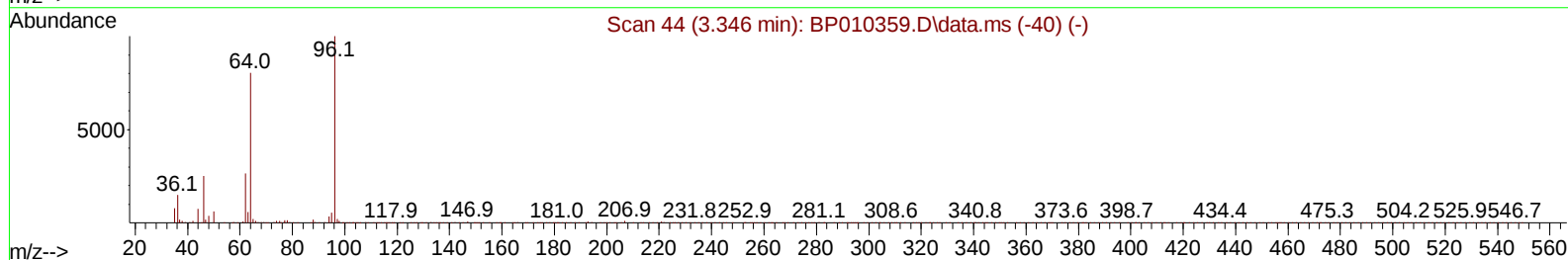
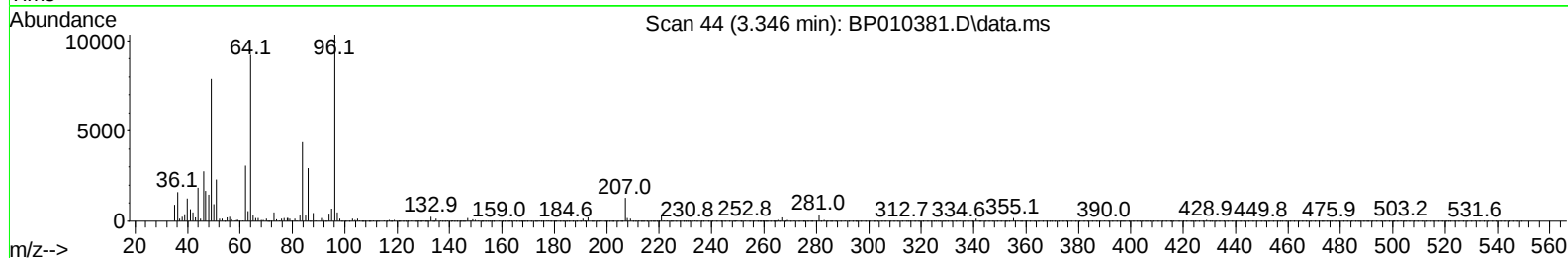
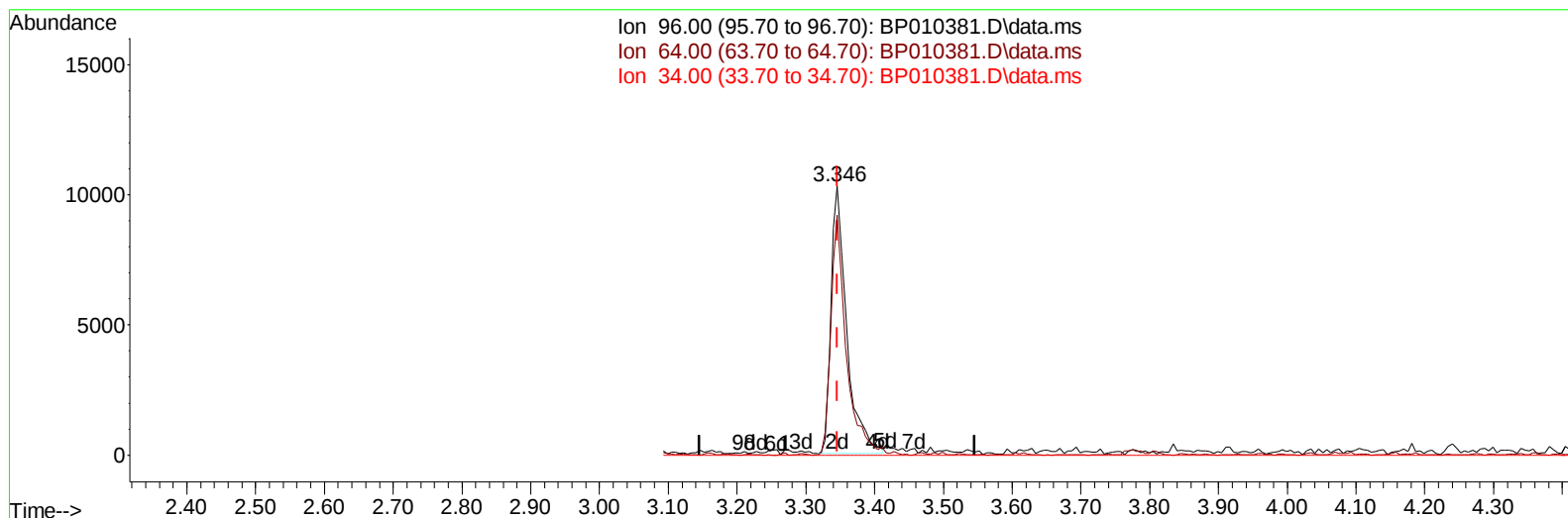
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TIC: BP010381.D\data.ms

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

3.346min (-0.000) 4.31 ng/uL m

response 16410

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	89.32#
34.00	0.00	0.00
0.00	0.00	0.00

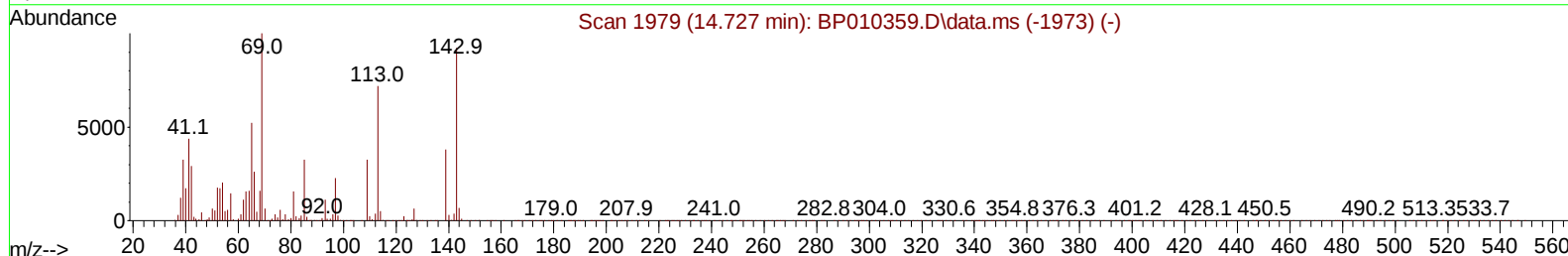
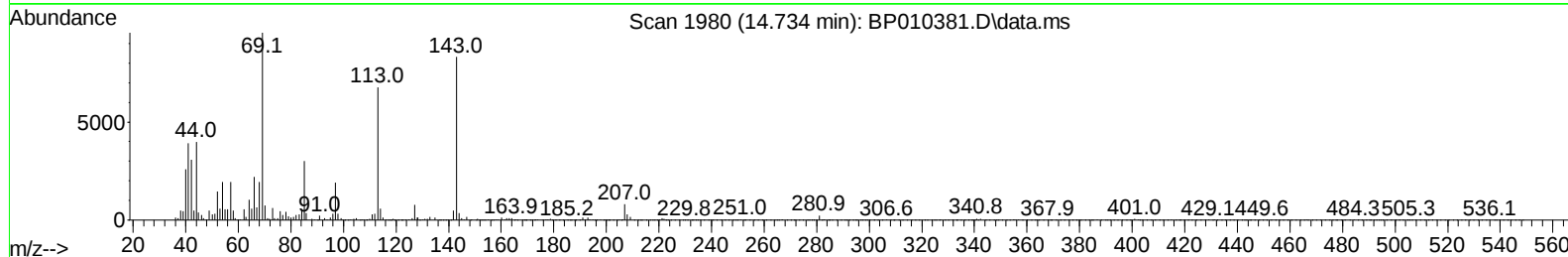
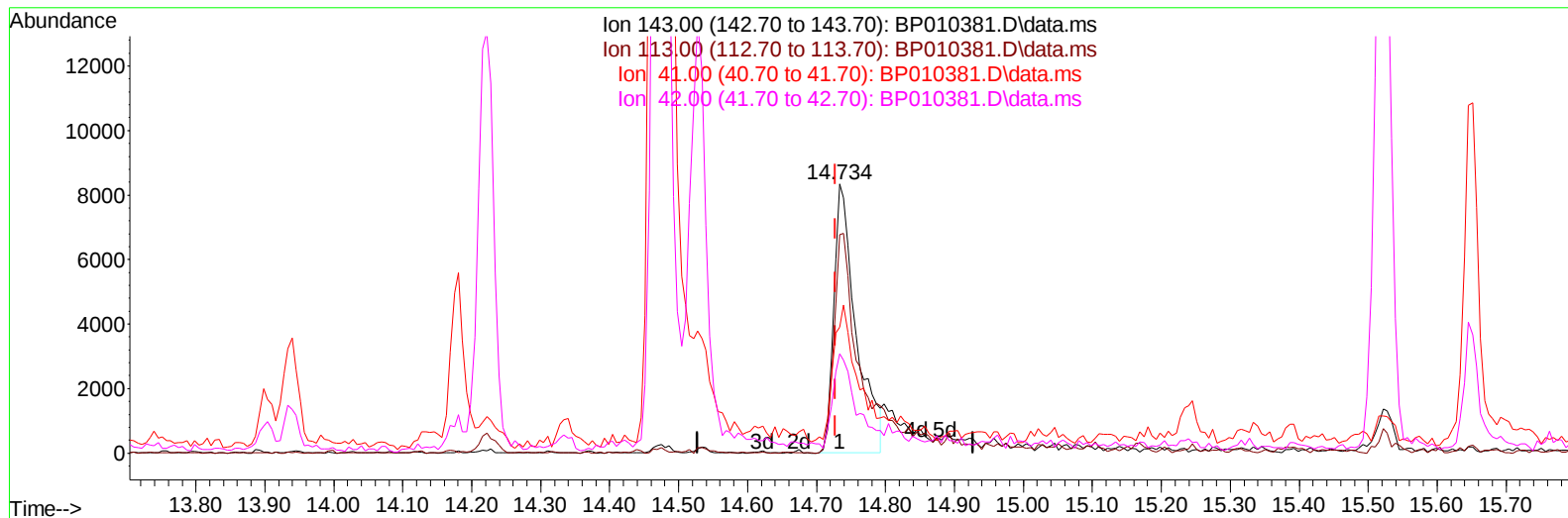
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Instrument :
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ClientSampleId :
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TIC: BP010381.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (+ 0.006) 3.02 ng/u1

response 18699

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	81.18#
41.00	23.20	46.86#
42.00	18.00	36.93#

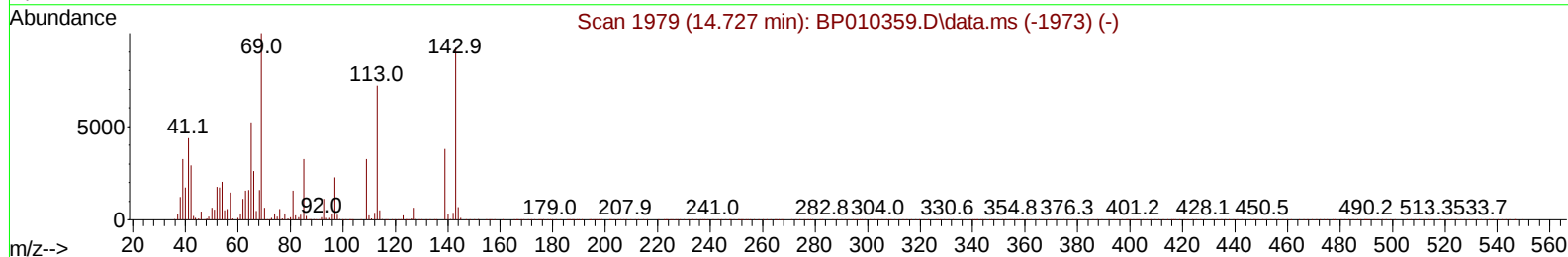
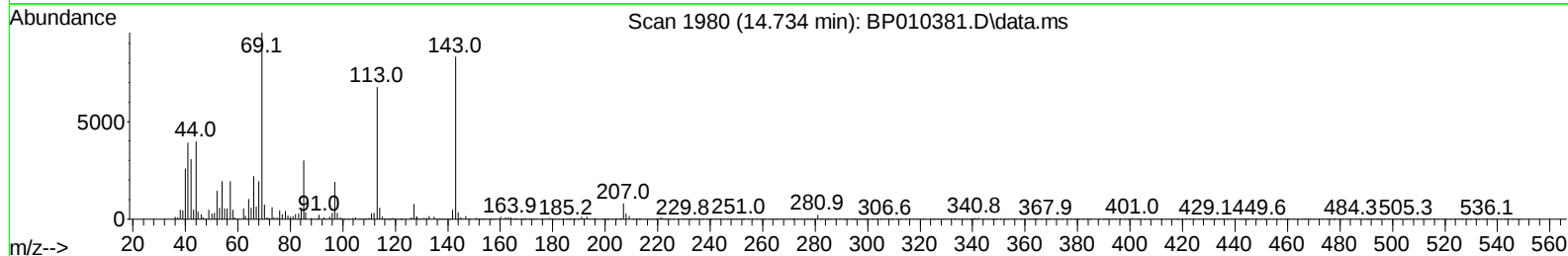
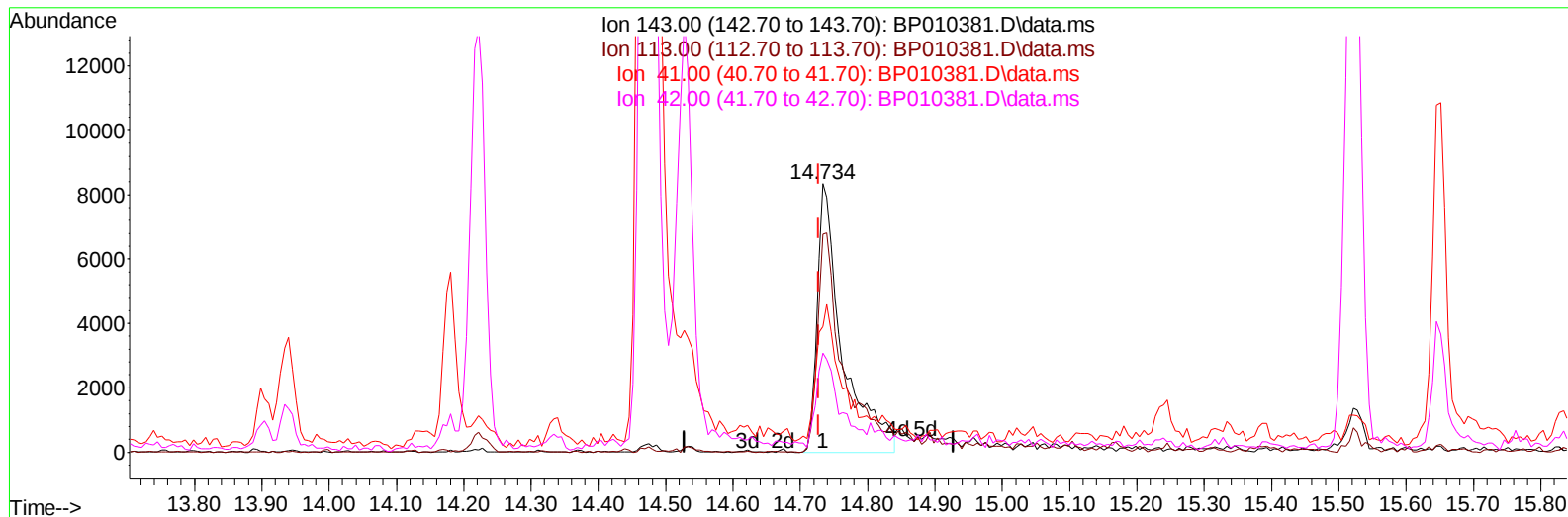
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ALS Vial : 29 Sample Multiplier: 1

Instrument :
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TIC: BP010381.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (+ 0.006) 3.49 ng/ul m

response 21648

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	81.18#
41.00	23.20	46.86#
42.00	18.00	36.93#

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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	7.899	152	162464	20.000	ng/ul	0.00
20) Naphthalene-d8	10.698	136	719910	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.528	164	472442	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.280	188	1059937	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.363	240	1046841	20.000	ng/ul	# 0.00
88) Perylene-d12	23.792	264	918854	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	16410m	4.305	ng/uL	0.00
4) Pyridine-d5	3.769	84	95688	8.842	ng/ul	0.00
7) Phenol-d5	7.063	99	83471	6.121	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	305367	33.331	ng/ul	0.00
11) 2-Chlorophenol-d4	7.428	132	260245	24.942	ng/ul	0.00
15) 4-Methylphenol-d8	8.604	113	180235	15.553	ng/ul	0.00
21) Nitrobenzene-d5	9.057	128	191247	34.325	ng/ul	0.00
24) 2-Nitrophenol-d4	9.781	143	210189	35.597	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.322	165	341442	30.582	ng/ul	0.00
31) 4-Chloroaniline-d4	10.834	131	378520	22.802	ng/ul	0.00
46) Dimethylphthalate-d6	13.939	166	1283618	36.877	ng/ul	0.00
49) Acenaphthylene-d8	14.222	160	1439218	35.846	ng/ul	0.00
54) 4-Nitrophenol-d4	14.734	143	21648m	3.494	ng/ul	0.00
60) Fluorene-d10	15.522	176	1085486	35.845	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	152600	24.617	ng/ul	0.00
73) Anthracene-d10	17.380	188	1769620	36.870	ng/ul	0.00
81) Pyrene-d10	19.616	212	2117320	38.085	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.633	264	1722071	37.323	ng/ul	0.00

Target Compounds Qvalue

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

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