

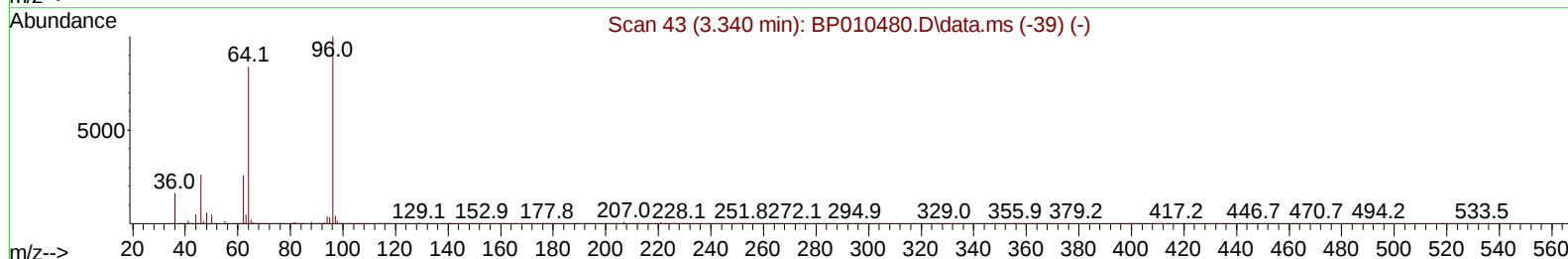
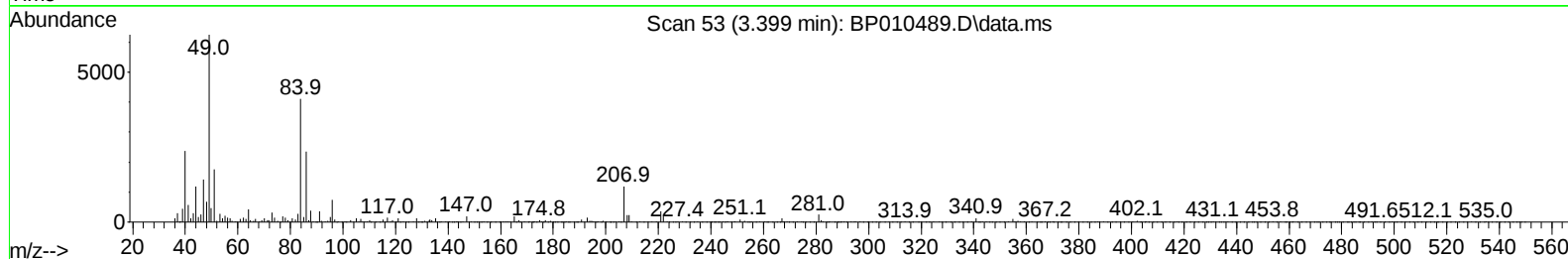
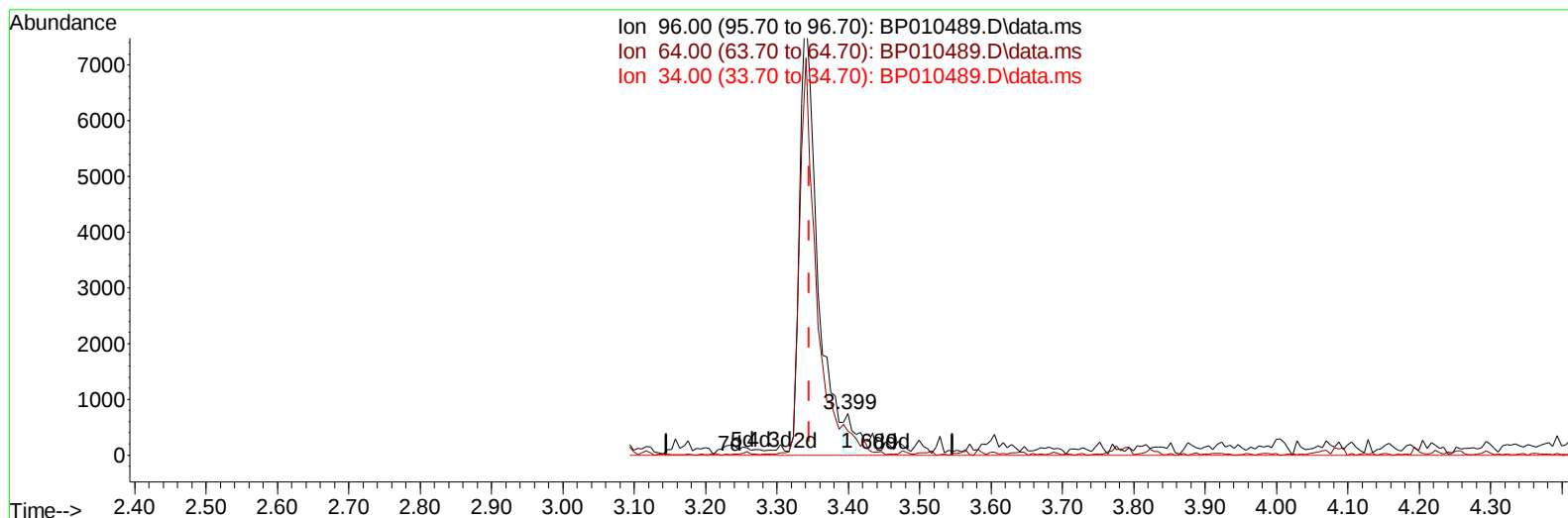
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052322\
 Data File : BP010489.D
 Acq On : 23 May 2022 22:16
 Operator : CG/JU
 Sample : N2955-02
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AA1

Manual IntegrationsAPPROVED

Quant Time: May 23 23:21:32 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 05/24/2022
 Supervised By :Yogesh Patel 05/26/2022



TIC: BP010489.D\data.ms

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

3.399min (+ 0.053) 0.16 ng/uL

response 503

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	57.62
34.00	0.00	0.00
0.00	0.00	0.00

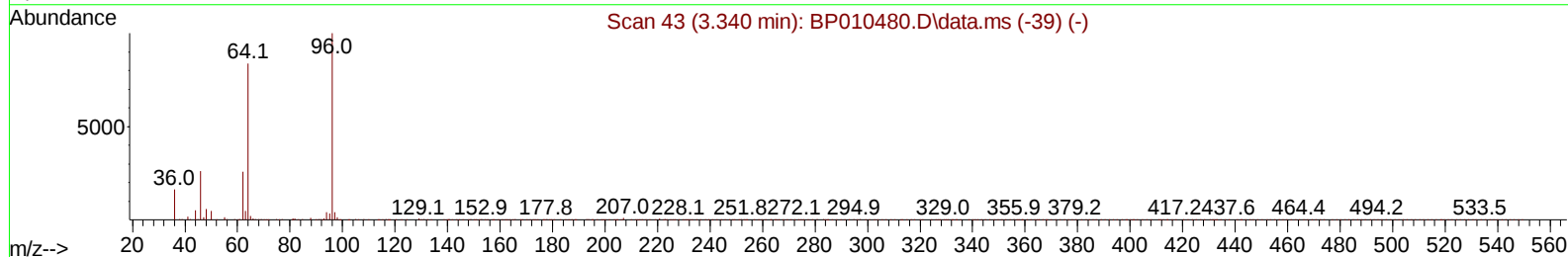
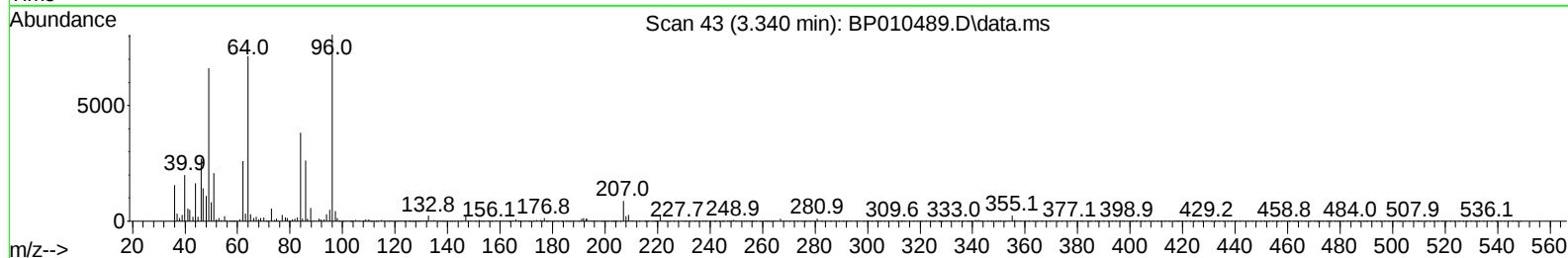
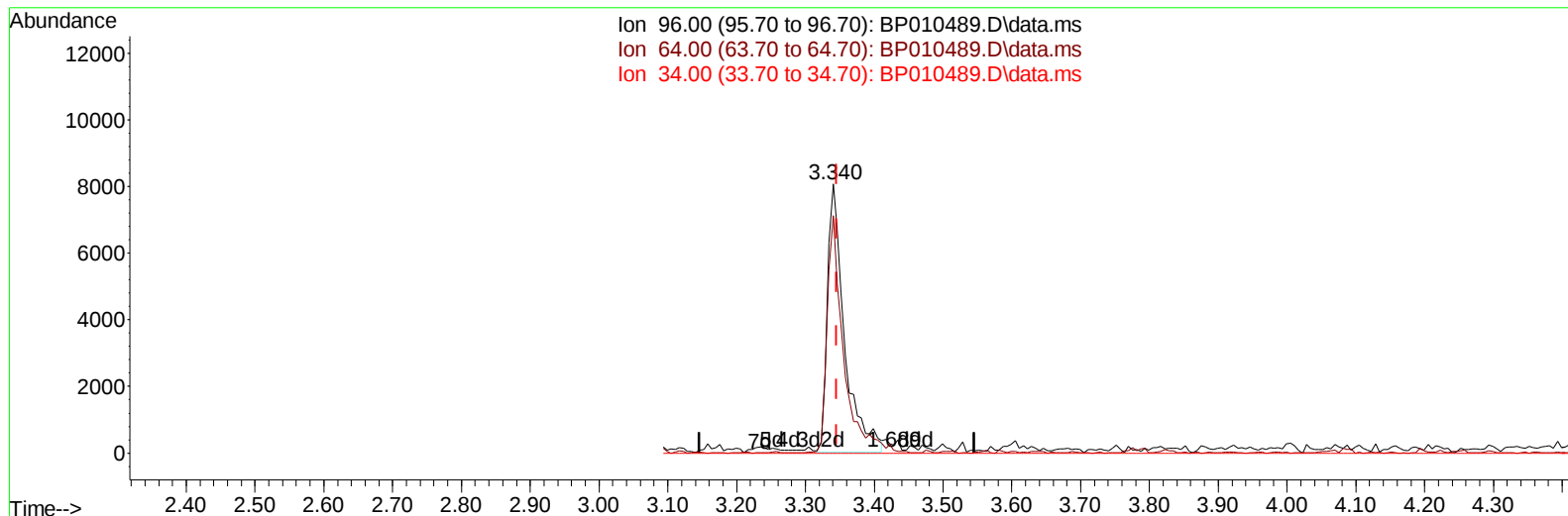
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TIC: BP010489.D\data.ms

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

3.340min (-0.006) 4.54 ng/uL m

response 13932

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	88.39#
34.00	0.00	0.00
0.00	0.00	0.00

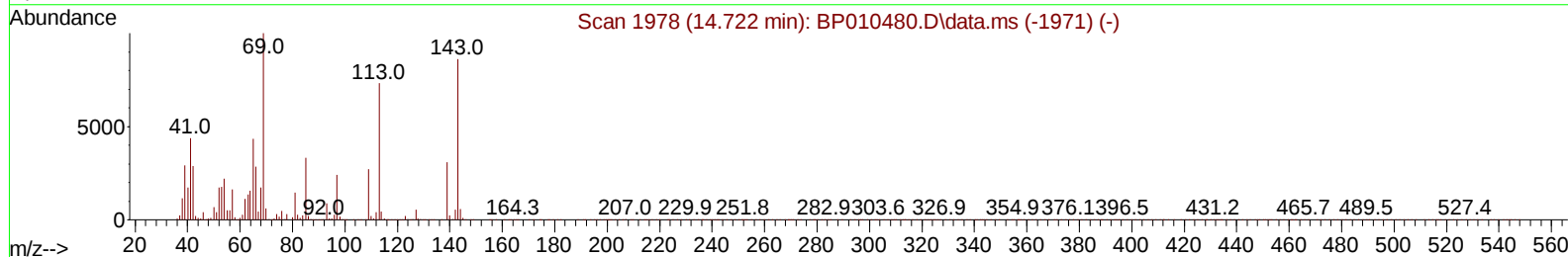
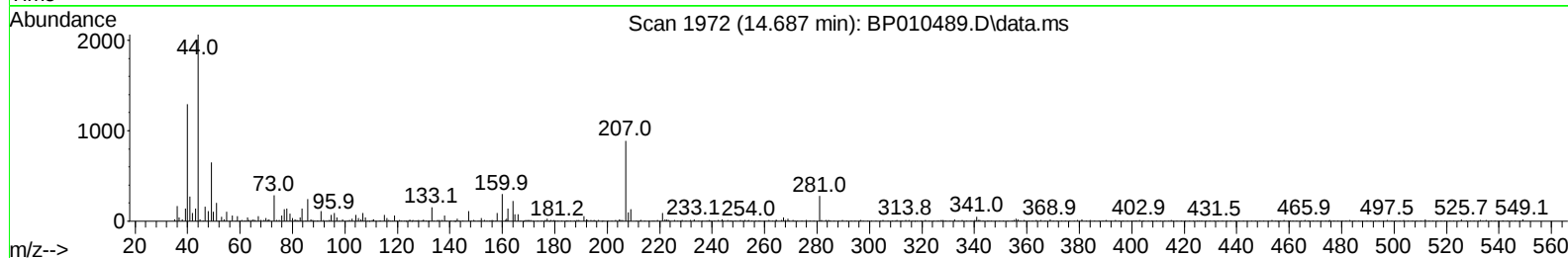
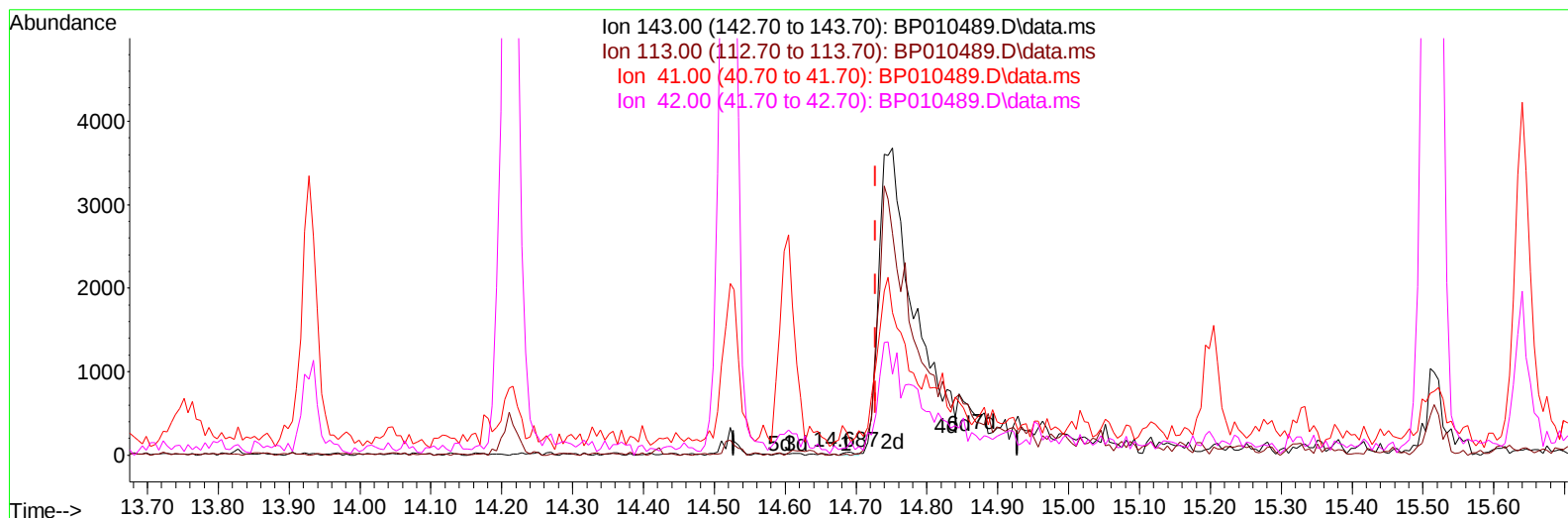
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ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: BP010489.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.687min (-0.041) 0.01 ng/u1

response 47

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	38.46#
41.00	23.20	700.00#
42.00	18.00	225.64#

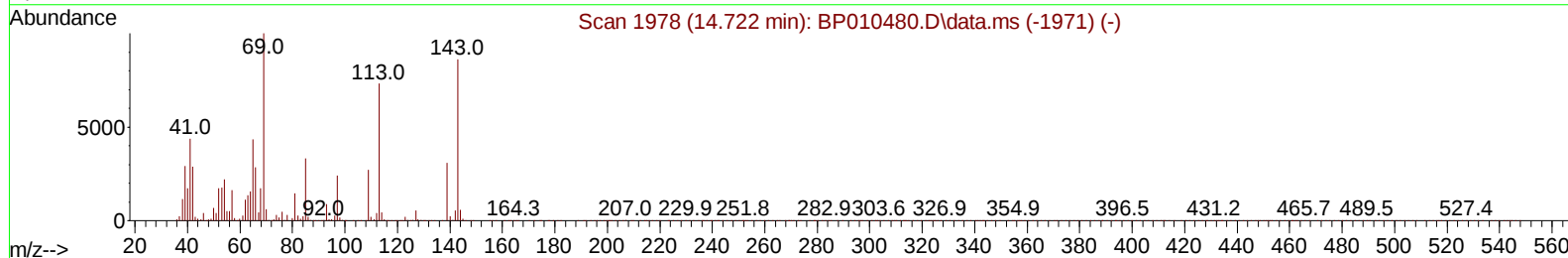
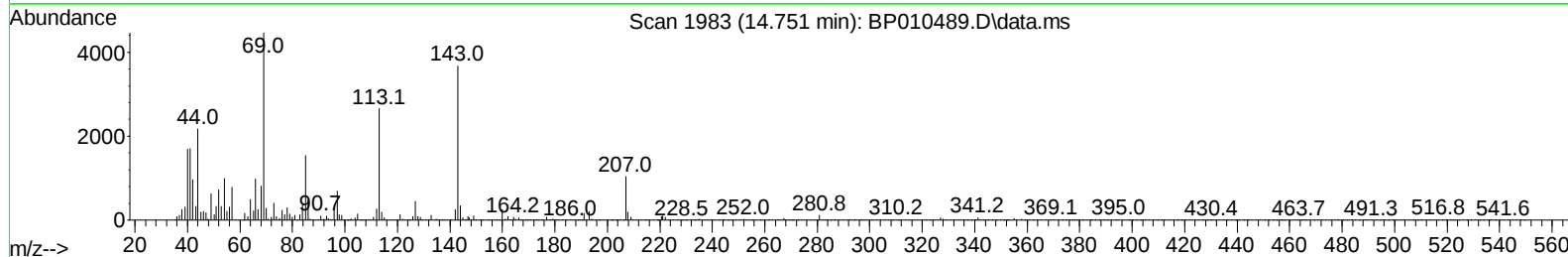
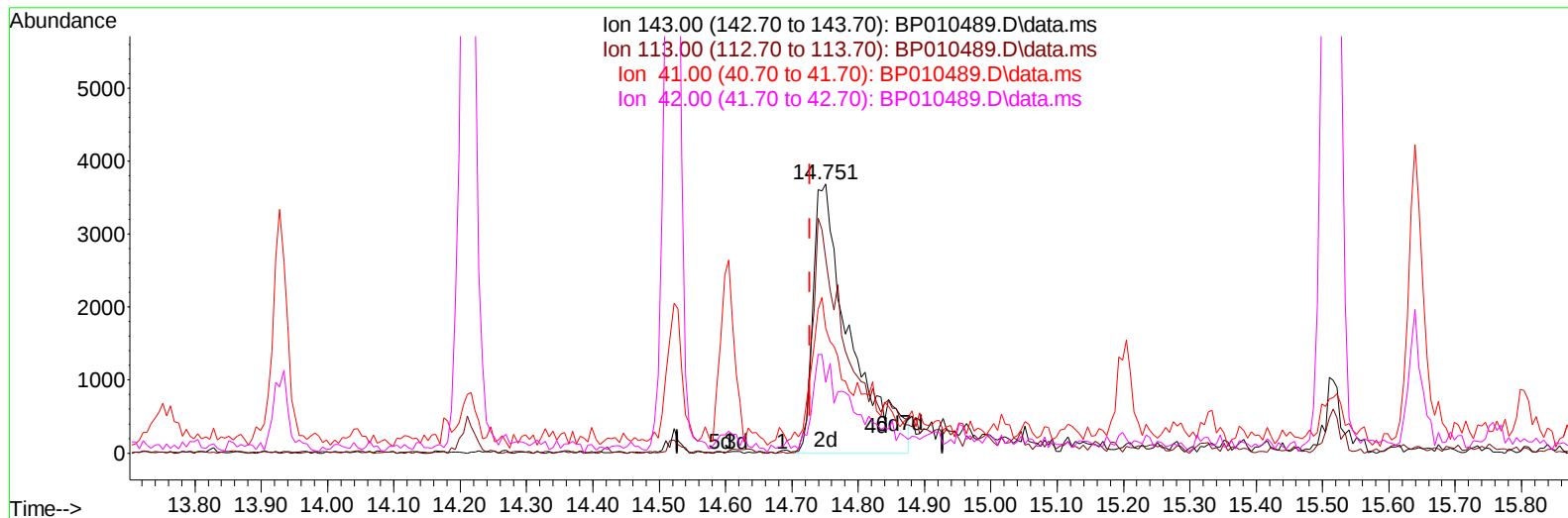
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Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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TIC: BP010489.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (+ 0.024) 2.59 ng/ul m

response 14226

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	72.16#
41.00	23.20	46.32#
42.00	18.00	26.27#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	130757	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.687	136	601568	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.522	164	419402	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.269	188	907796	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.351	240	848929	20.000	ng/ul	#-0.01
88) Perylene-d12	23.774	264	688025	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	13932m	4.542	ng/uL	0.00
4) Pyridine-d5	3.776	84	39043	4.482	ng/ul	0.01
7) Phenol-d5	7.058	99	64389	5.867	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	225418	30.571	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.422	132	200035	23.820	ng/ul	-0.01
15) 4-Methylphenol-d8	8.599	113	138116	14.809	ng/ul	-0.01
21) Nitrobenzene-d5	9.046	128	143315	30.782	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.769	143	133761	27.110	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.310	165	242691	26.013	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.828	131	289308	20.856	ng/ul	-0.01
46) Dimethylphthalate-d6	13.928	166	974258	31.529	ng/ul	-0.01
49) Acenaphthylene-d8	14.210	160	1104007	30.974	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.751	143	14226m	2.587	ng/ul	0.02
60) Fluorene-d10	15.516	176	850396	31.634	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	58132	10.949	ng/ul	0.00
73) Anthracene-d10	17.369	188	1451657	35.315	ng/ul	-0.01
81) Pyrene-d10	19.604	212	1744610	38.697	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.616	264	1250879	36.207	ng/ul	-0.02

Target Compounds Qvalue

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

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