

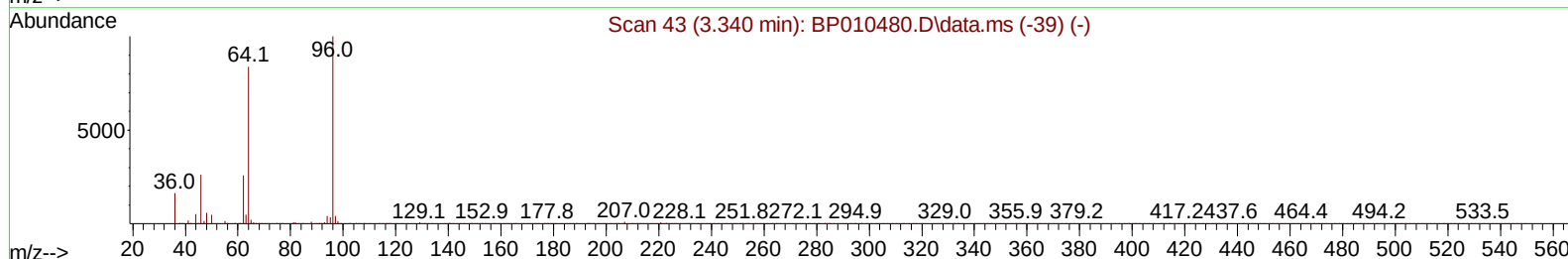
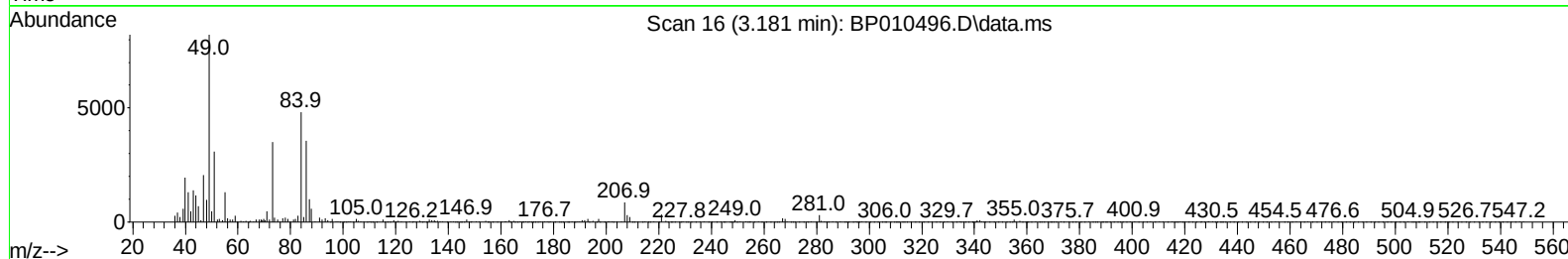
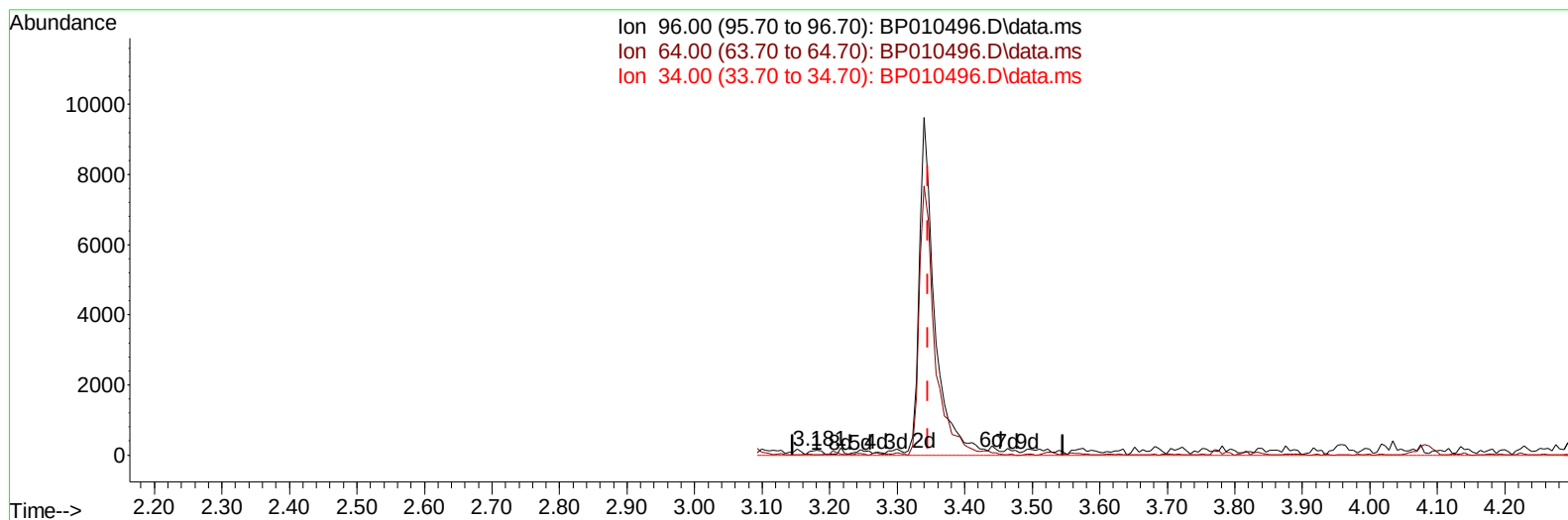
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052322\
 Data File : BP010496.D
 Acq On : 24 May 2022 02:13
 Operator : CG/JU
 Sample : N2950-02
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBTE7

Manual IntegrationsAPPROVED

Quant Time: May 24 03:08:45 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: BP010496.D\data.ms

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

3.181min (-0.165) 0.04 ng/uL

response 146

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	54.55
34.00	0.00	0.00
0.00	0.00	0.00

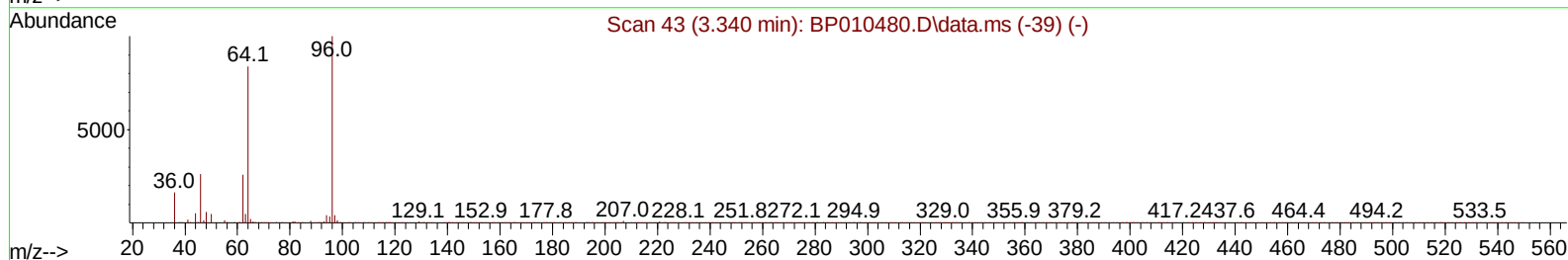
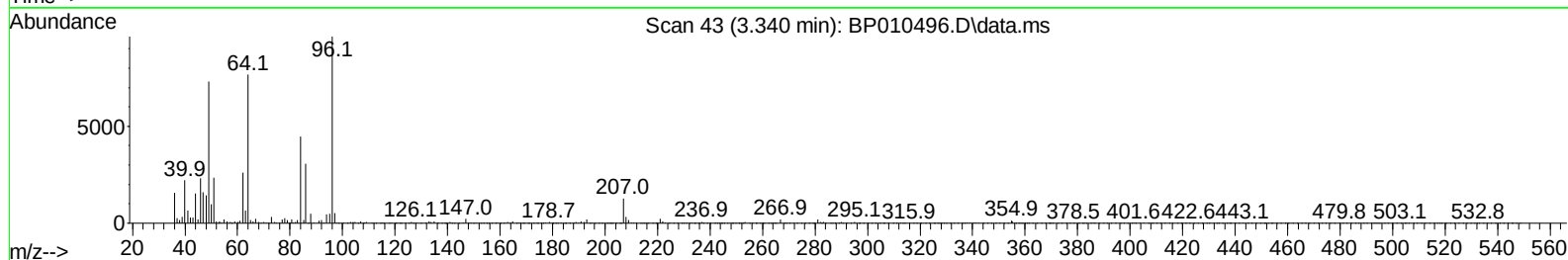
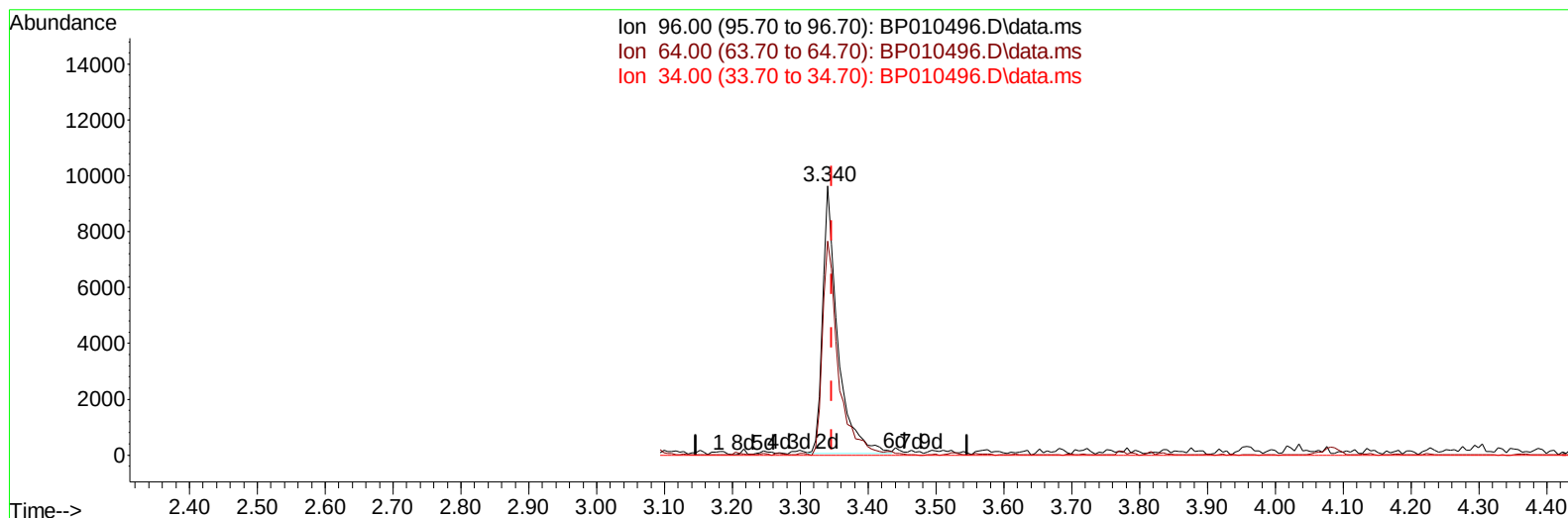
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TIC: BP010496.D\data.ms

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

3.340min (-0.006) 4.17 ng/uL m

response 14867

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	79.63#
34.00	0.00	0.00
0.00	0.00	0.00

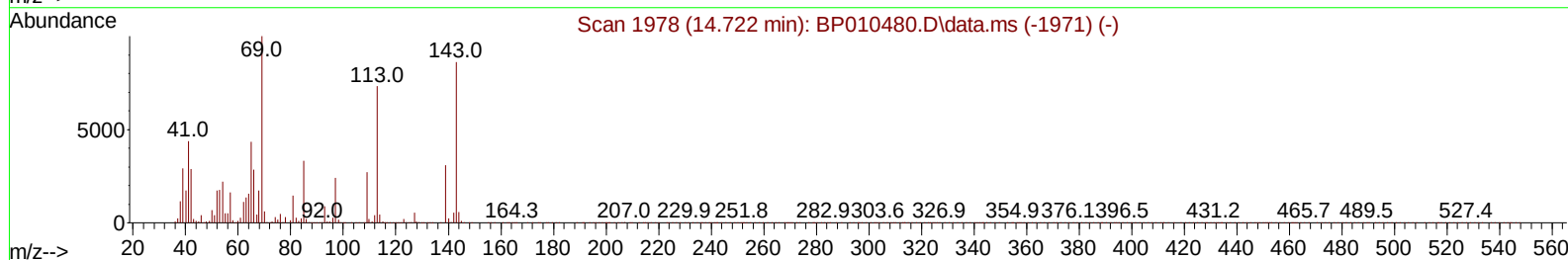
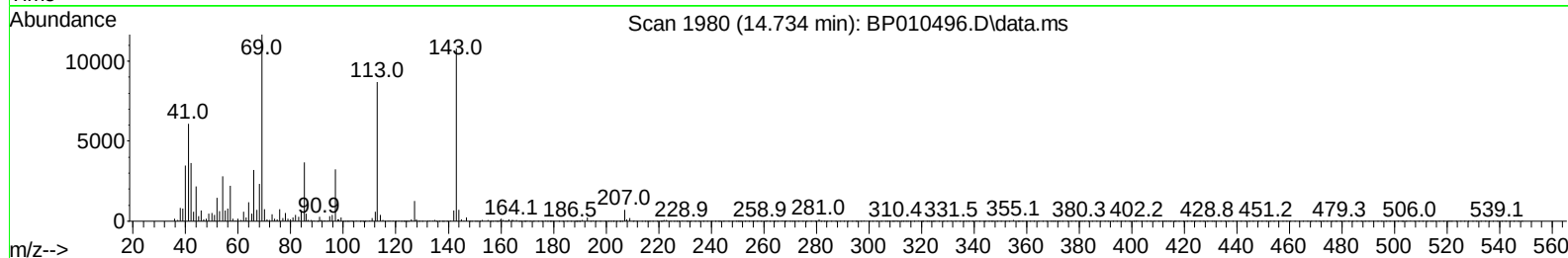
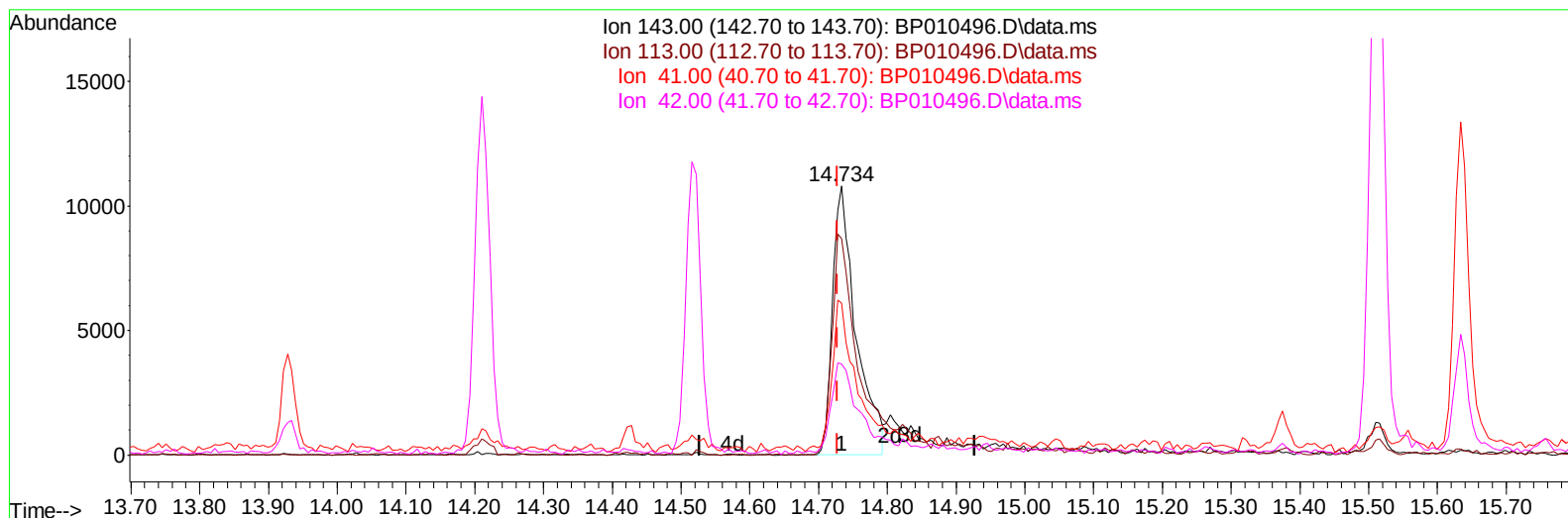
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052322\
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 Operator : CG/JU
 Sample : N2950-02
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
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TIC: BP010496.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (+ 0.006) 4.23 ng/ul

response 25543

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	80.54#
41.00	23.20	56.41#
42.00	18.00	33.76#

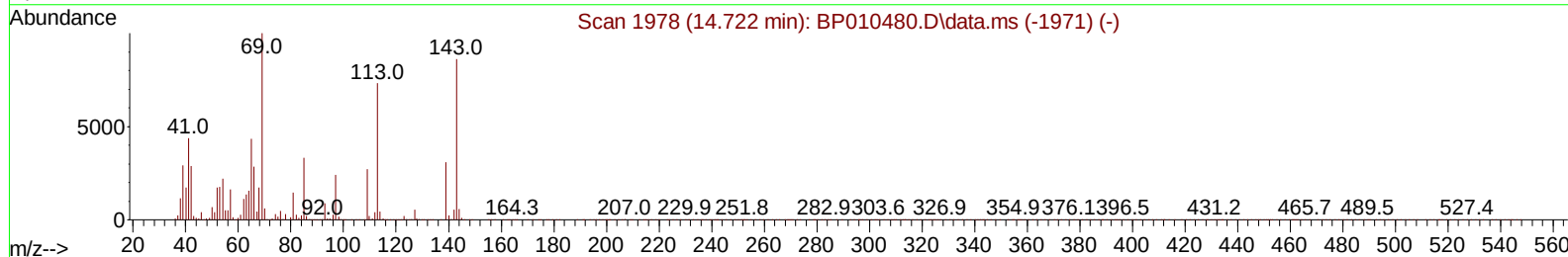
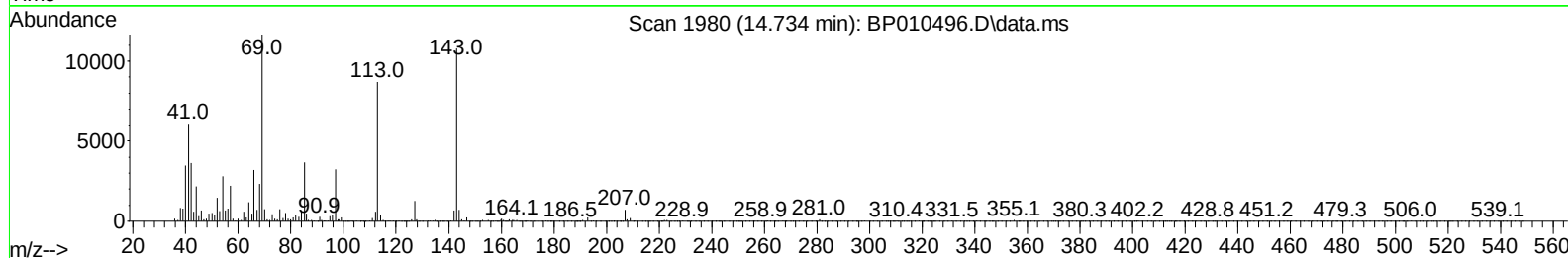
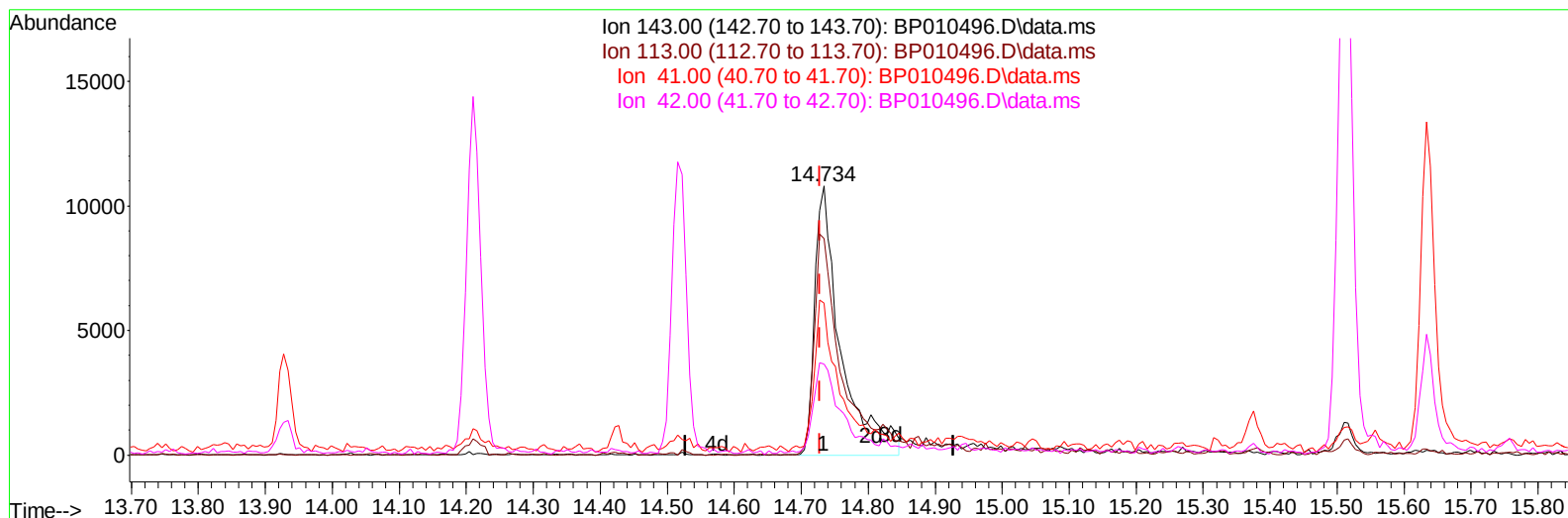
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 Sample : N2950-02
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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TIC: BP010496.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.734min (+ 0.006) 4.82 ng/ul m

response 29096

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	80.54#
41.00	23.20	56.41#
42.00	18.00	33.76#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	151792	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.687	136	685357	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.516	164	460607	20.000	ng/ul	-0.02
64) Phenanthrene-d10	17.269	188	984694	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.351	240	796947	20.000	ng/ul	-0.01
88) Perylene-d12	23.774	264	635078	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	14867m	4.175	ng/uL	0.00
4) Pyridine-d5	3.775	84	41089	4.064	ng/ul	0.01
7) Phenol-d5	7.058	99	89171	6.999	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	292914	34.220	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.422	132	255646	26.223	ng/ul	-0.01
15) 4-Methylphenol-d8	8.599	113	175546	16.213	ng/ul	-0.01
21) Nitrobenzene-d5	9.046	128	184997	34.877	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.769	143	189121	33.644	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.310	165	310040	29.170	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.828	131	292732	18.523	ng/ul	-0.01
46) Dimethylphthalate-d6	13.928	166	1250235	36.841	ng/ul	-0.01
49) Acenaphthylene-d8	14.210	160	1383433	35.342	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.734	143	29096m	4.817	ng/ul	0.00
60) Fluorene-d10	15.516	176	1081291	36.624	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.634	200	182975	31.773	ng/ul	-0.01
73) Anthracene-d10	17.375	188	1856085	41.627	ng/ul	0.00
81) Pyrene-d10	19.604	212	2153962	50.893	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.616	264	1371856	43.019	ng/ul	-0.02

Target Compounds Qvalue

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

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