

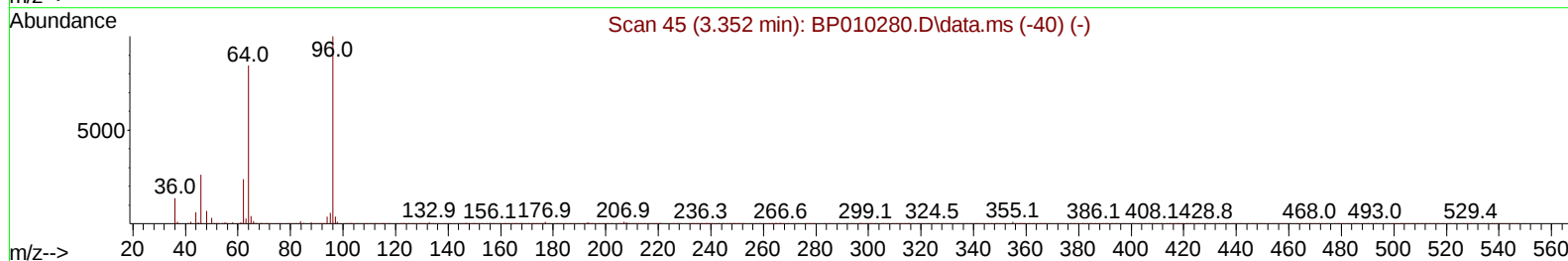
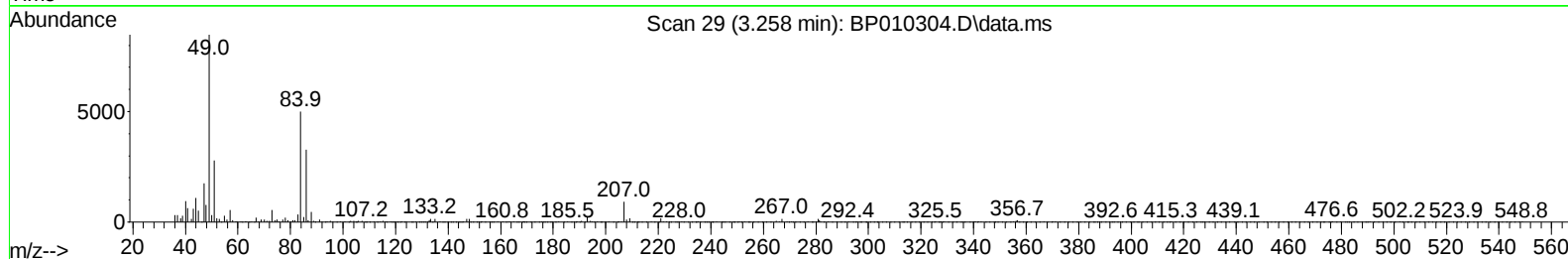
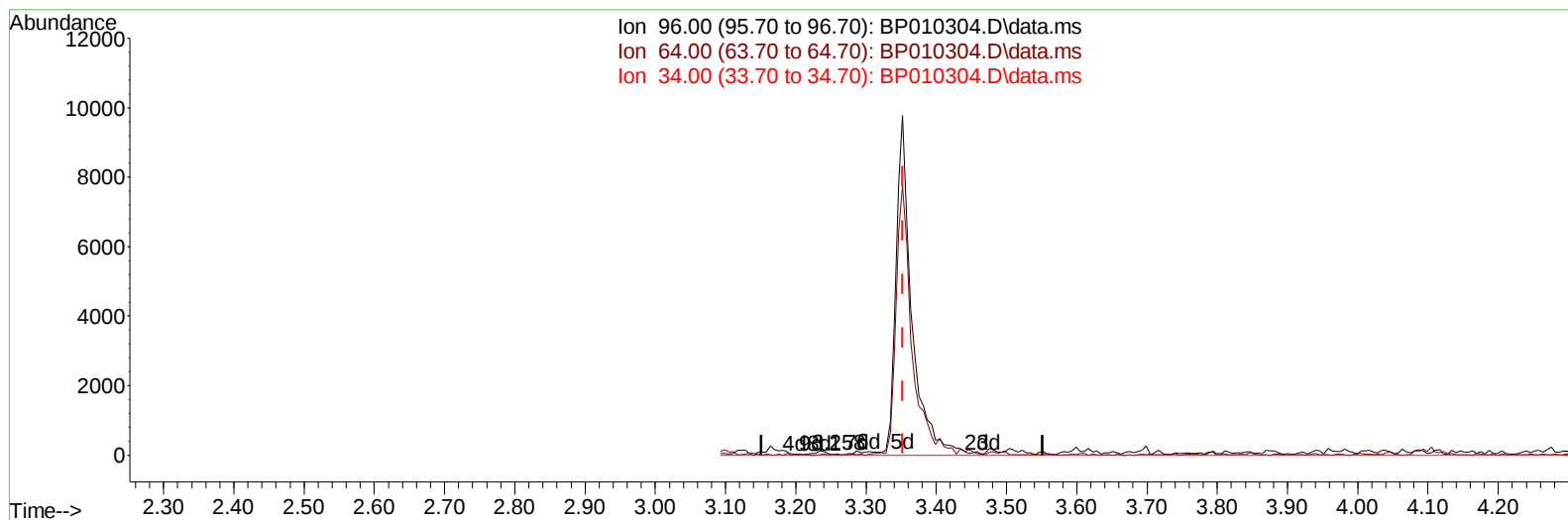
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
 Data File : BP010304.D
 Acq On : 12 May 2022 02:09
 Operator : CG/JU
 Sample : N2707-07
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXL88

Manual IntegrationsAPPROVED

Quant Time: May 12 02:59:46 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu May 12 01:10:15 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/12/2022
 Supervised By :Yogesh Patel 05/17/2022



TIC: BP010304.D\data.ms

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

3.258min (-0.094) 0.01 ng/uL

response 21

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	56.00
34.00	0.00	0.00
0.00	0.00	0.00

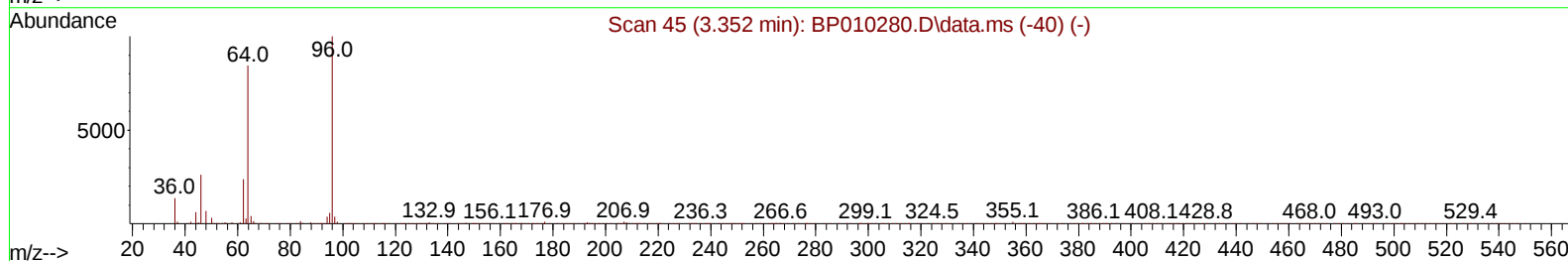
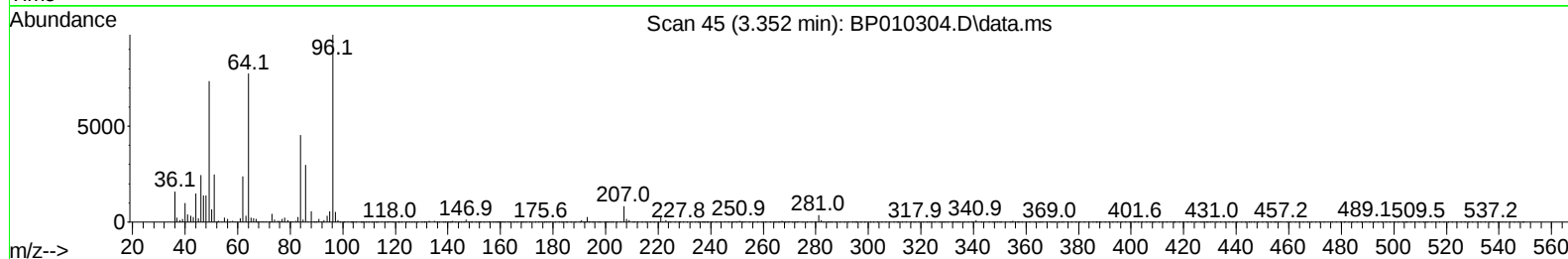
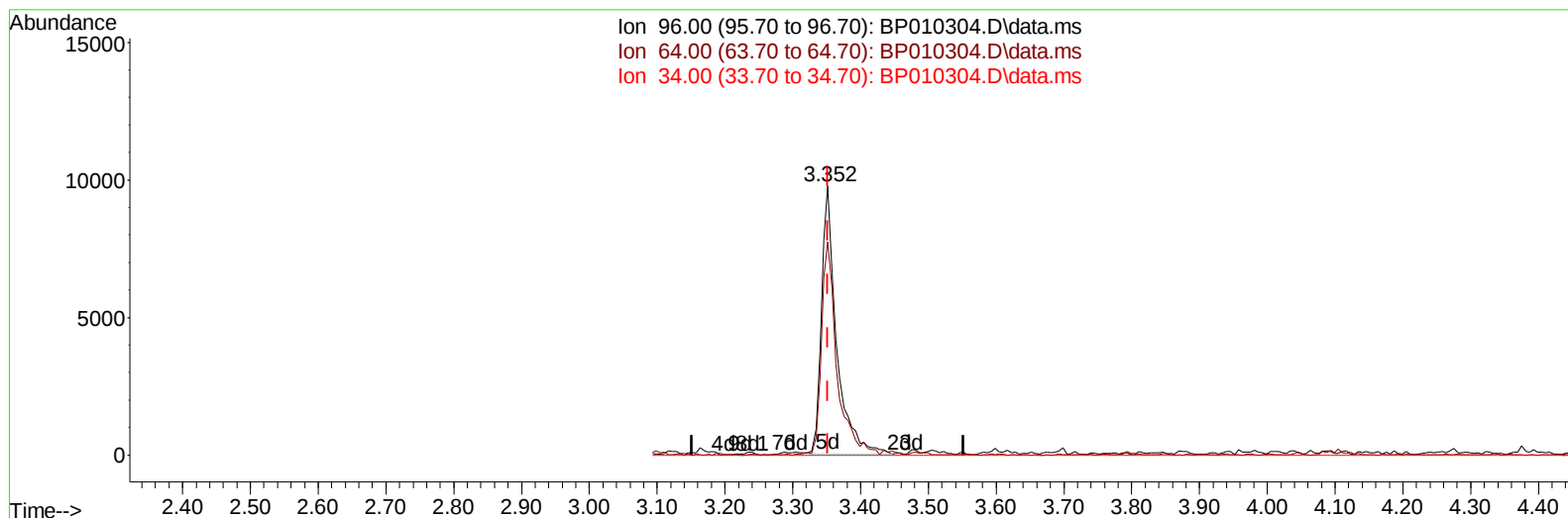
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TIC: BP010304.D\data.ms

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

3.352min (+ 0.000) 4.36 ng/uL m

response 15043

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	79.23#
34.00	0.00	0.00
0.00	0.00	0.00

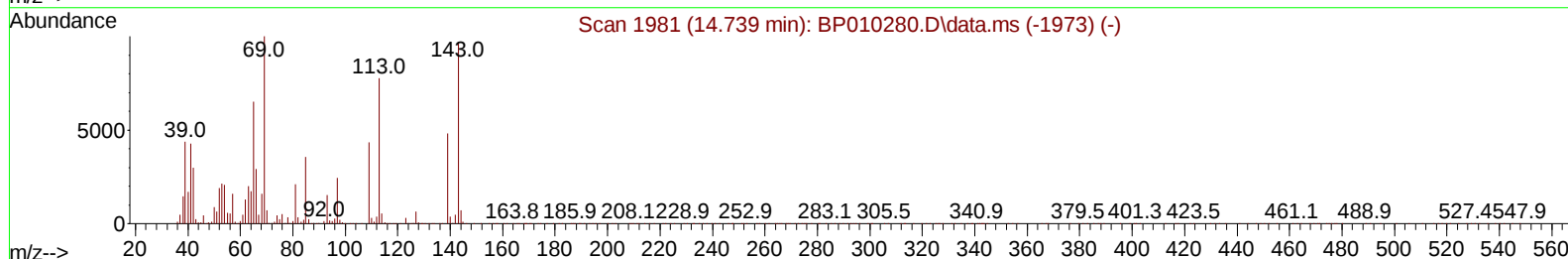
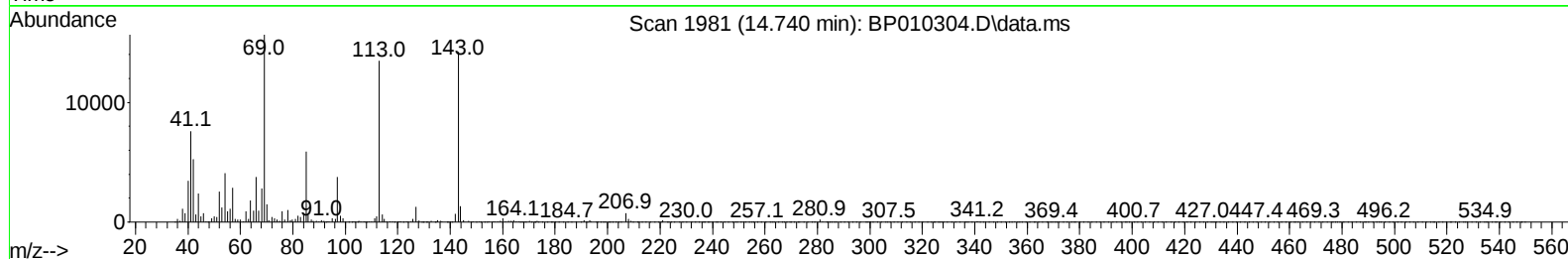
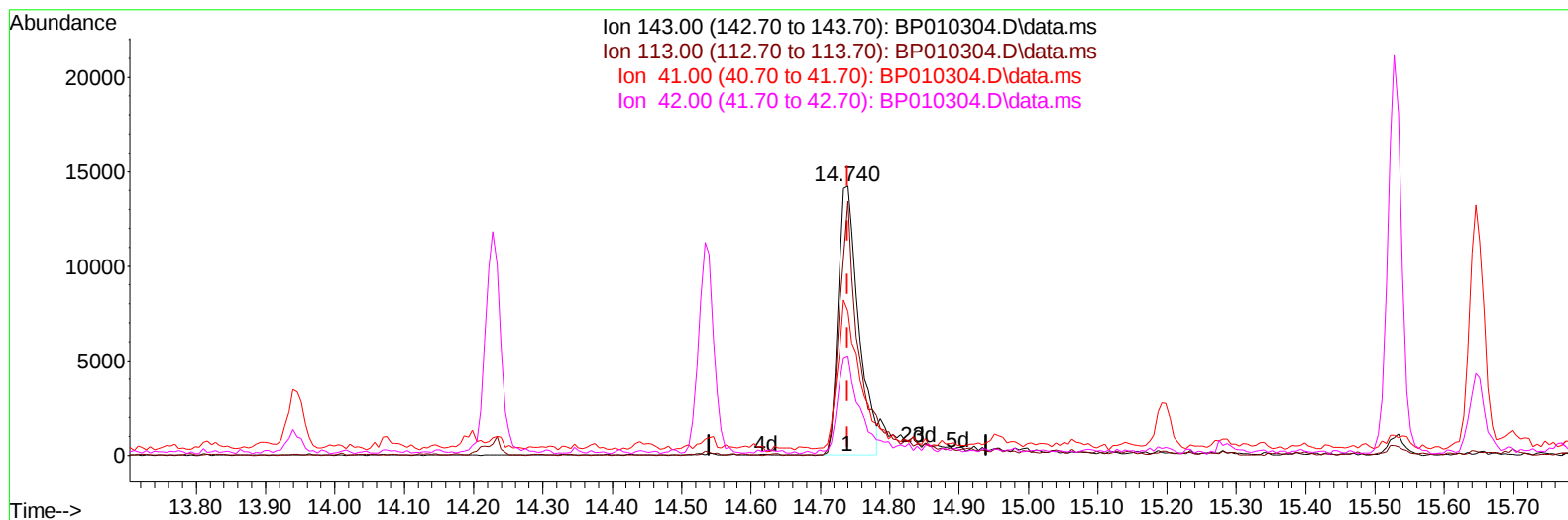
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Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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TIC: BP010304.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.740min (+ 0.000) 4.23 ng/u1

response 28769

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	94.52#
41.00	23.20	53.37#
42.00	18.00	36.99#

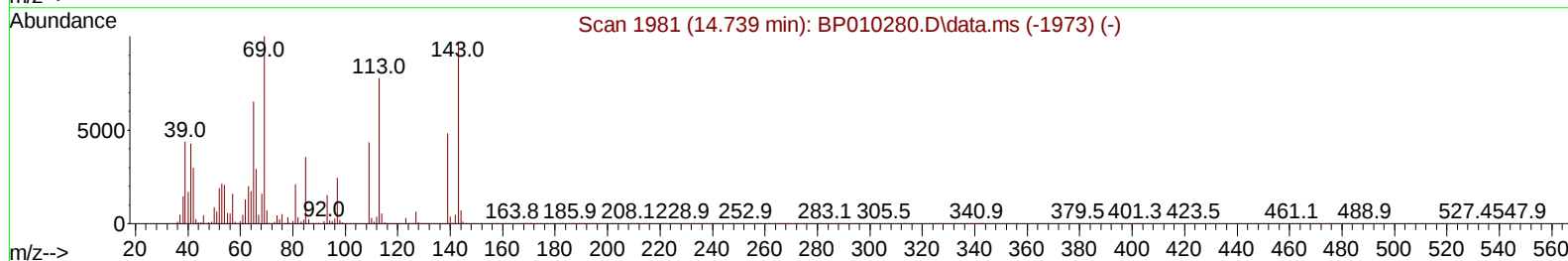
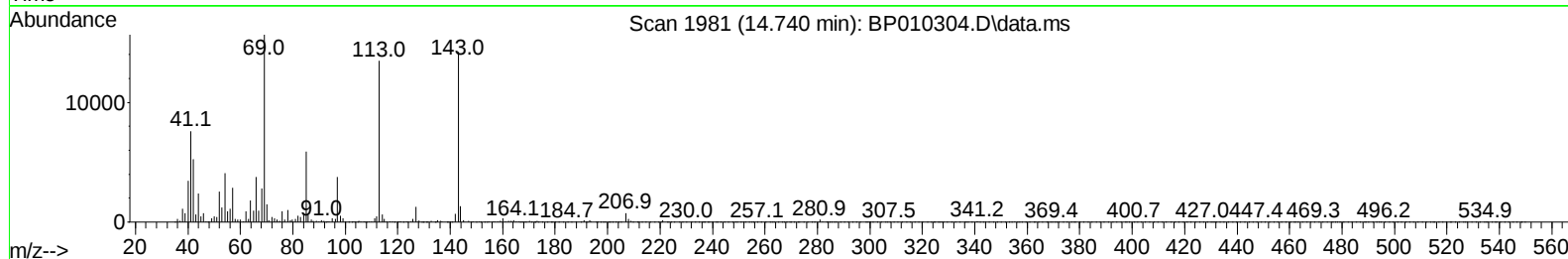
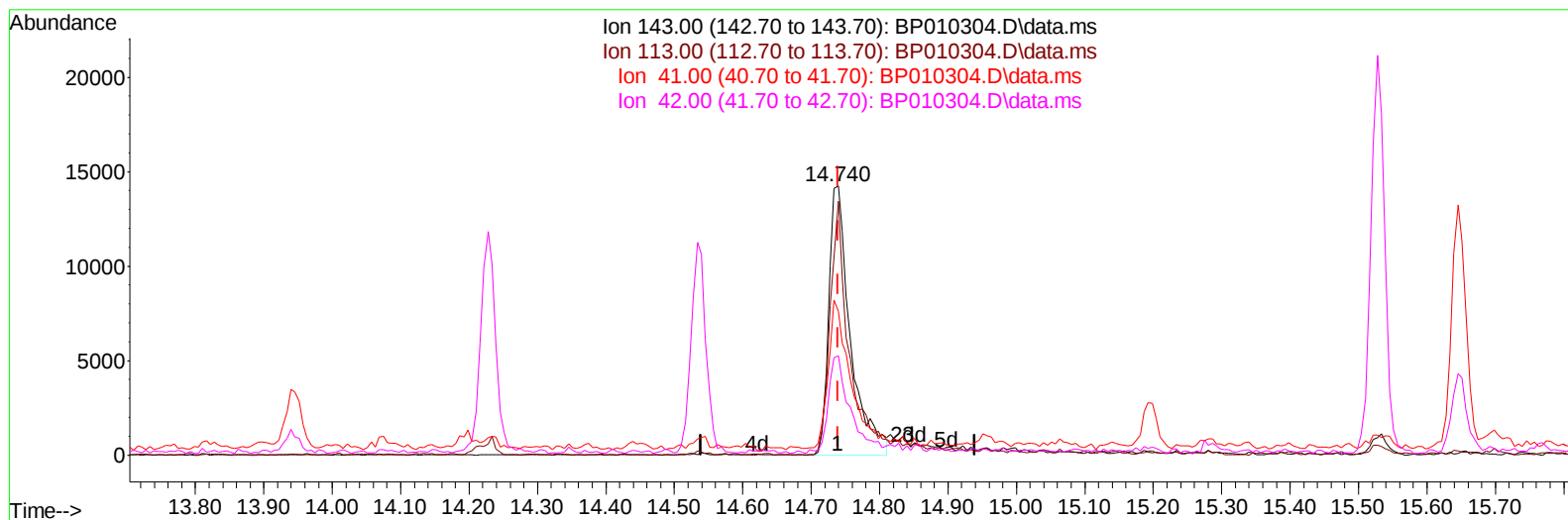
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Instrument :
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TIC: BP010304.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.740min (+ 0.000) 4.57 ng/ul m

response 31107

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	94.52#
41.00	23.20	53.37#
42.00	18.00	36.99#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.905	152	146501	20.000	ng/ul	0.00
20) Naphthalene-d8	10.704	136	628244	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.540	164	445343	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.286	188	993627	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.369	240	1014738	20.000	ng/ul	# 0.00
88) Perylene-d12	23.798	264	994064	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	15043m	4.362	ng/uL	0.00
4) Pyridine-d5	3.781	84	37912	3.951	ng/ul	0.01
7) Phenol-d5	7.069	99	72760	5.347	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	251608	29.891	ng/ul	0.00
11) 2-Chlorophenol-d4	7.440	132	217077	21.662	ng/ul	0.00
15) 4-Methylphenol-d8	8.610	113	147341	13.171	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128	153003	30.374	ng/ul	0.00
24) 2-Nitrophenol-d4	9.787	143	160091	27.874	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165	268992	25.732	ng/ul	0.00
31) 4-Chloroaniline-d4	10.846	131	269691	17.475	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	1119274	31.548	ng/ul	0.00
49) Acenaphthylene-d8	14.228	160	1230757	29.096	ng/ul	0.00
54) 4-Nitrophenol-d4	14.740	143	31107m	4.571	ng/ul	0.00
60) Fluorene-d10	15.528	176	967556	32.307	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	182495	26.821	ng/ul	0.00
73) Anthracene-d10	17.386	188	1676490	34.454	ng/ul	0.00
81) Pyrene-d10	19.616	212	2024985	36.112	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.645	264	1884672	35.551	ng/ul	-0.01

Target Compounds Qvalue

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

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