

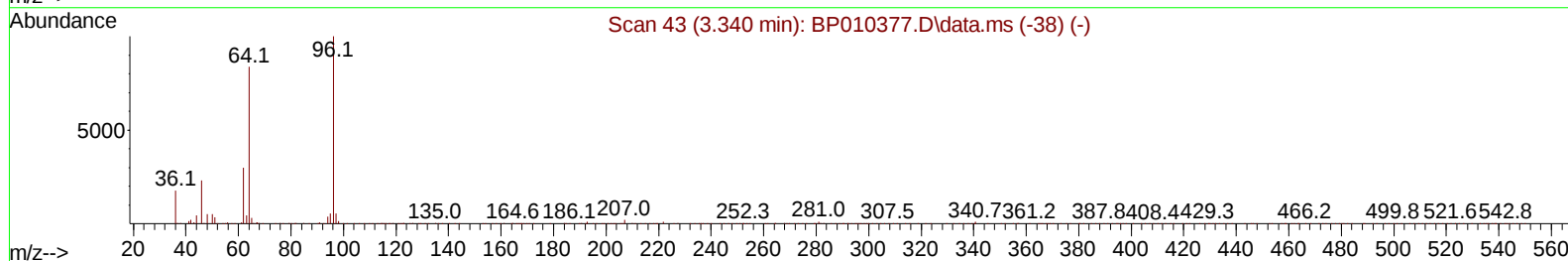
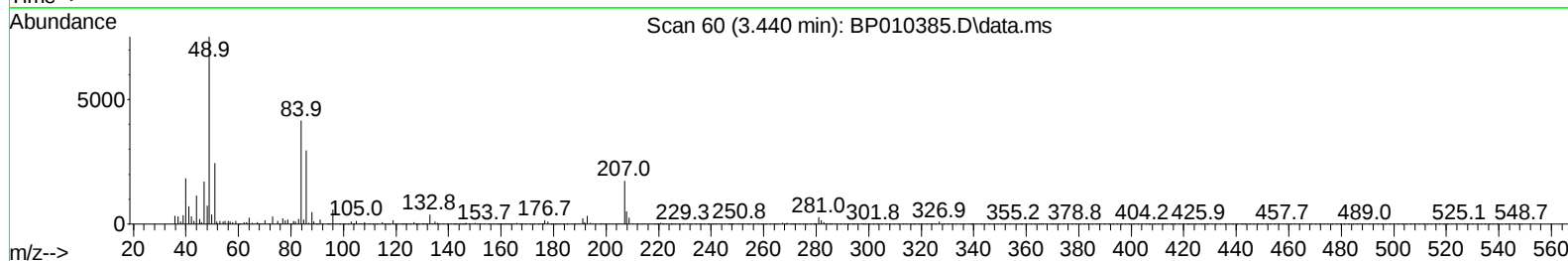
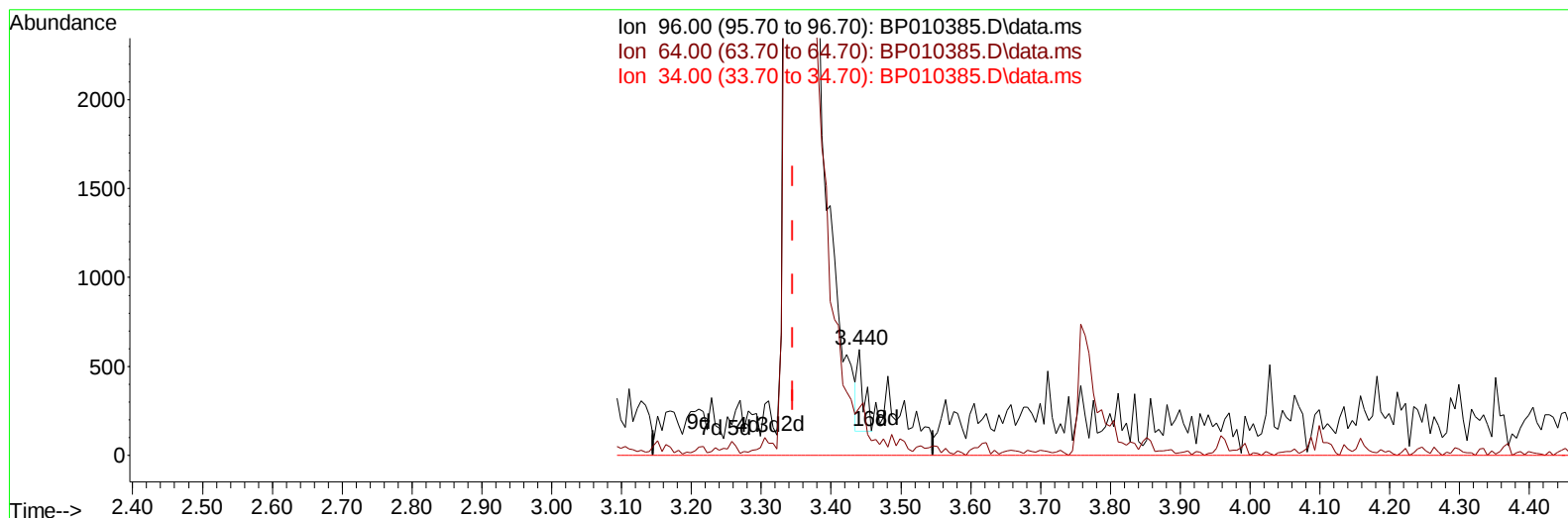
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051622\
 Data File : BP010385.D
 Acq On : 17 May 2022 09:02
 Operator : CG/JU
 Sample : PB144810BL
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK810

Manual IntegrationsAPPROVED

Quant Time: May 18 00:56:43 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 05/18/2022
 Supervised By :mohammad ahmed 05/18/2022



TIC: BP010385.D\data.ms

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

3.440min (+ 0.094) 0.07 ng/uL

response	286	
Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	44.95
34.00	0.00	0.00
0.00	0.00	0.00

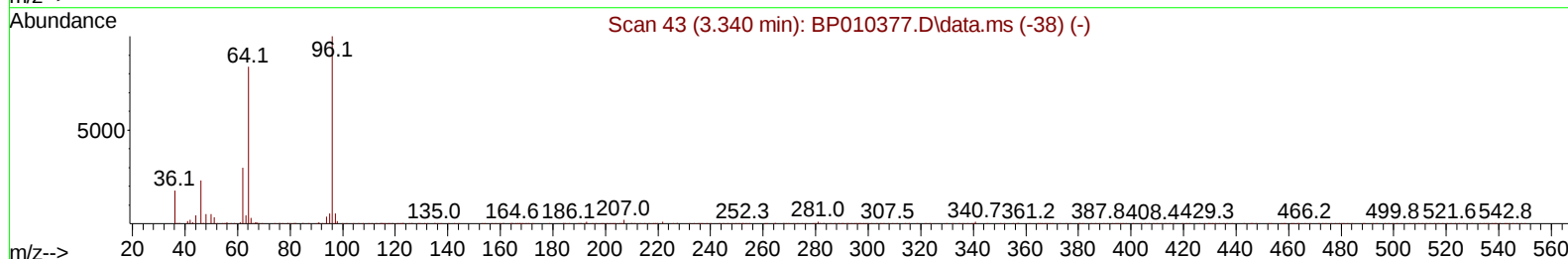
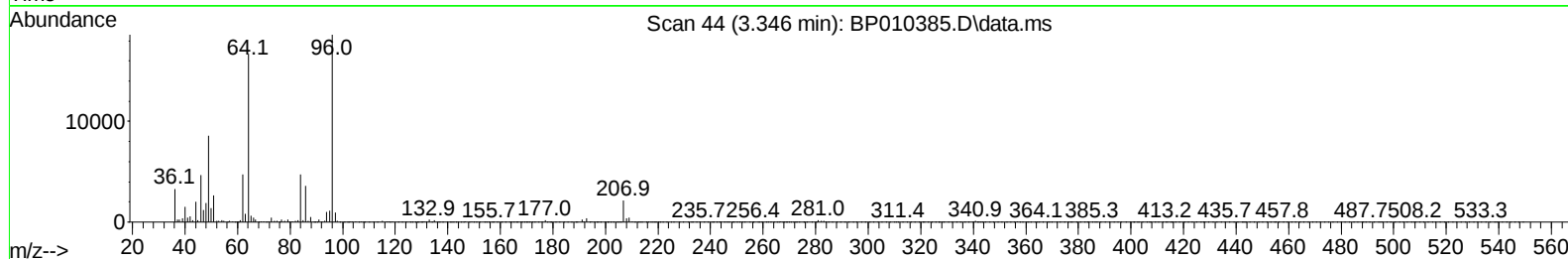
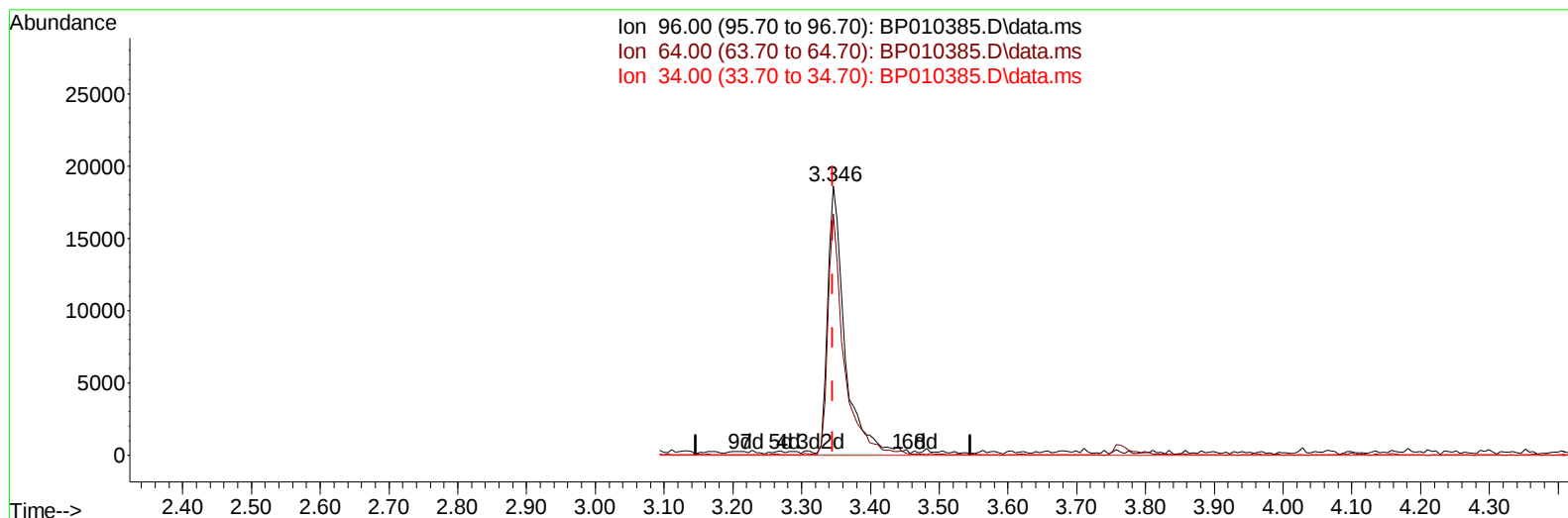
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TIC: BP010385.D\data.ms

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

3.346min (+ 0.000) 7.09 ng/uL m

response 31126

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	89.77#
34.00	0.00	0.00
0.00	0.00	0.00

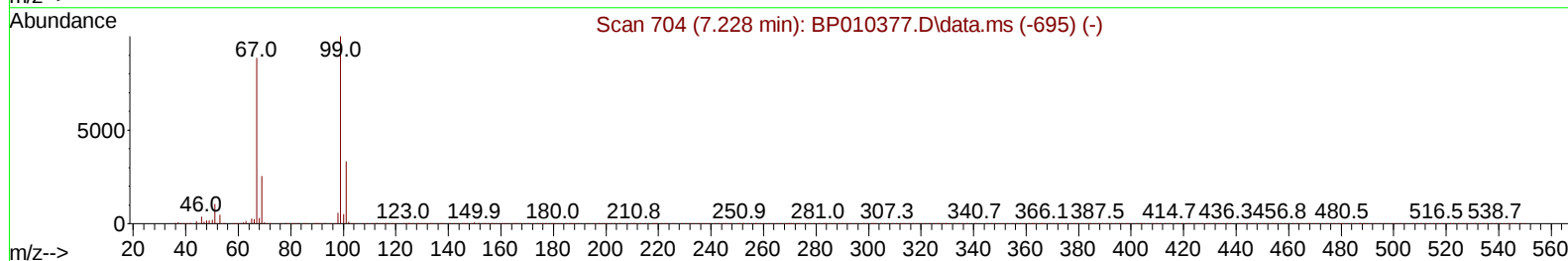
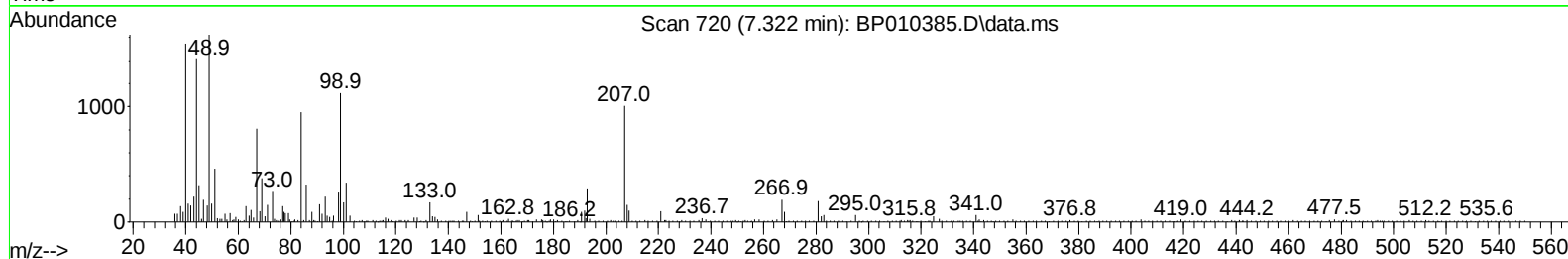
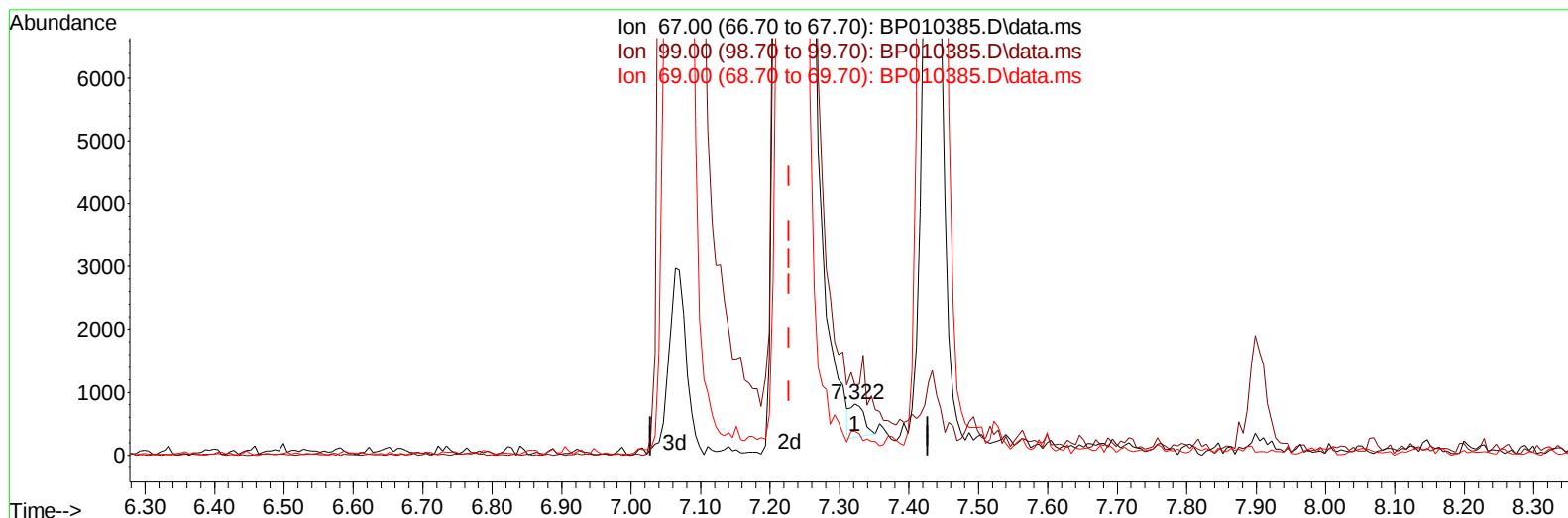
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 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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TIC: BP010385.D\data.ms

(9) Bis-(2-Chloroethyl)ether-d8 (S)

7.322min (+ 0.094) 0.07 ng/u1

response 745

Ion	Exp%	Act%
67.00	100.00	100.00
99.00	165.50	137.56
69.00	40.10	46.55
0.00	0.00	0.00

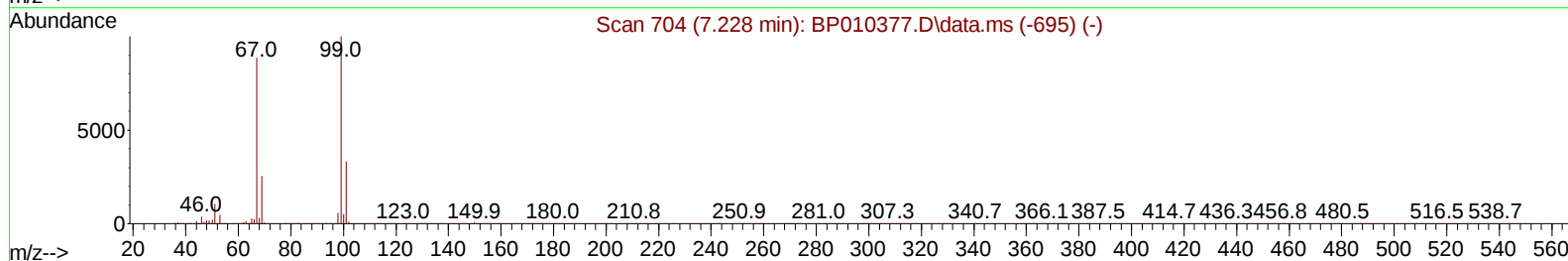
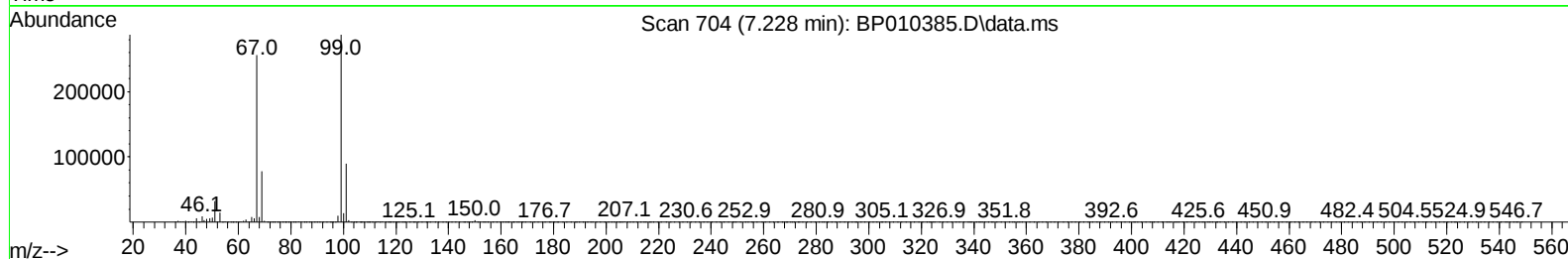
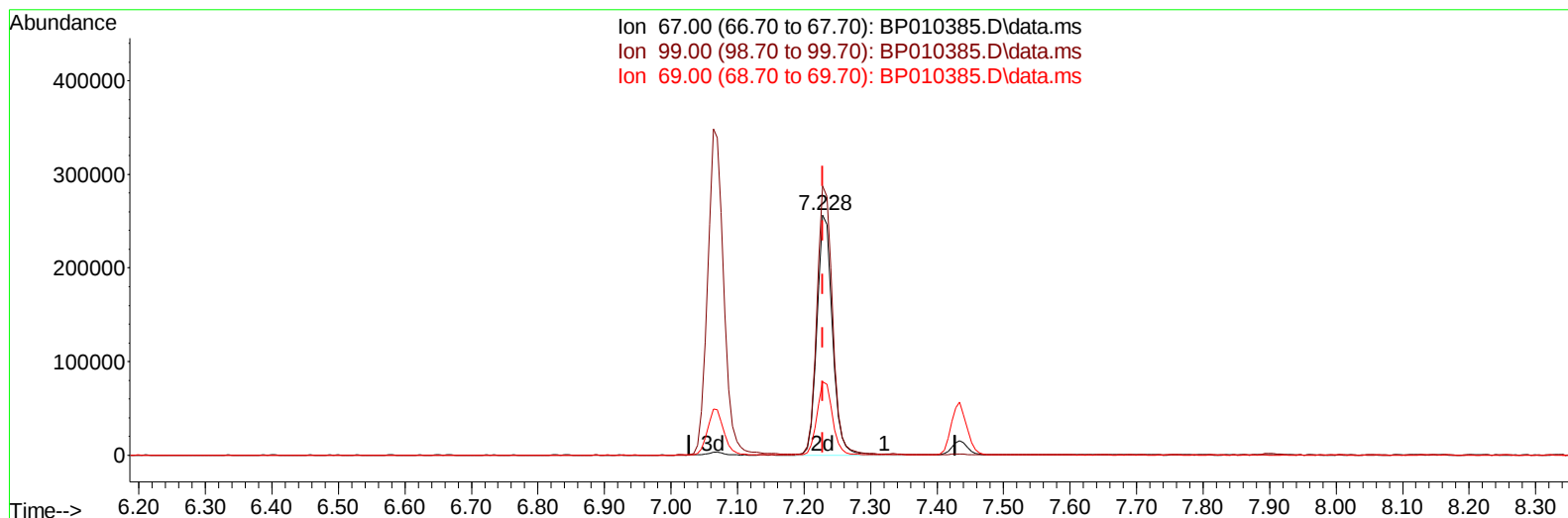
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TIC: BP010385.D\data.ms

(9) Bis-(2-Chloroethyl)ether-d8 (S)

7.228min (+ 0.000) 39.00 ng/ul m

response 411433

Ion	Exp%	Act%
67.00	100.00	100.00
99.00	165.50	111.99#
69.00	40.10	30.54#
0.00	0.00	0.00

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Instrument :
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Manual IntegrationsAPPROVED

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.905	152	187057	20.000	ng/ul	0.00
20) Naphthalene-d8	10.704	136	810263	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.540	164	549060	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.292	188	1187349	20.000	ng/ul	# 0.01
79) Chrysene-d12	21.375	240	1152740	20.000	ng/ul	# 0.01
88) Perylene-d12	23.810	264	1038162	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.346	96	31126m	7.093	ng/uL	0.00
4) Pyridine-d5	3.764	84	452476	36.313	ng/ul	0.00
7) Phenol-d5	7.063	99	589085	37.521	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	411433m	39.004	ng/ul	0.00
11) 2-Chlorophenol-d4	7.434	132	471771	39.270	ng/ul	0.00
15) 4-Methylphenol-d8	8.610	113	508462	38.108	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128	255130	40.685	ng/ul	0.00
24) 2-Nitrophenol-d4	9.787	143	275050	41.388	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165	475346	37.828	ng/ul	0.00
31) 4-Chloroaniline-d4	10.840	131	696398	37.272	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	1643351	40.624	ng/ul	0.00
49) Acenaphthylene-d8	14.228	160	1862676	39.919	ng/ul	0.00
54) 4-Nitrophenol-d4	14.734	143	258364	35.882	ng/ul	0.00
60) Fluorene-d10	15.534	176	1387991	39.439	ng/ul	0.01
65) 4,6-Dinitro-2-methylph...	15.651	200	215276	31.001	ng/ul	0.00
73) Anthracene-d10	17.392	188	2186035	40.659	ng/ul	0.01
81) Pyrene-d10	19.622	212	2626792	42.909	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.657	264	2207025	42.337	ng/ul	0.02

Target Compounds Qvalue

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

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