

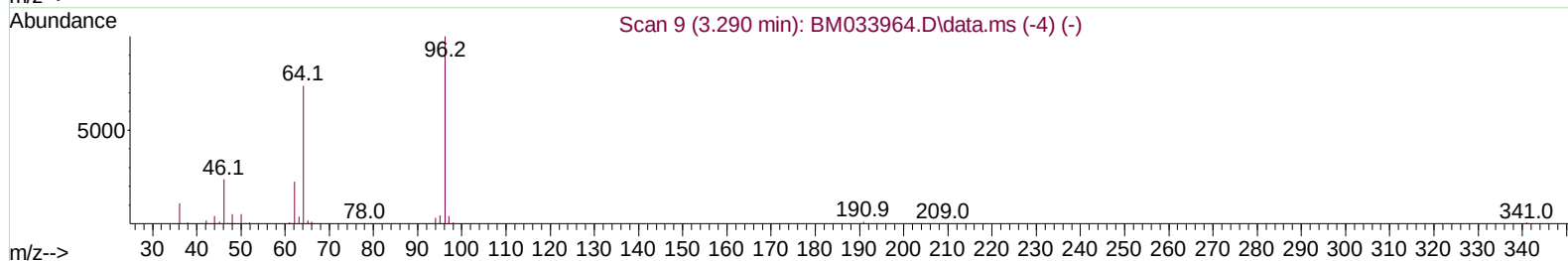
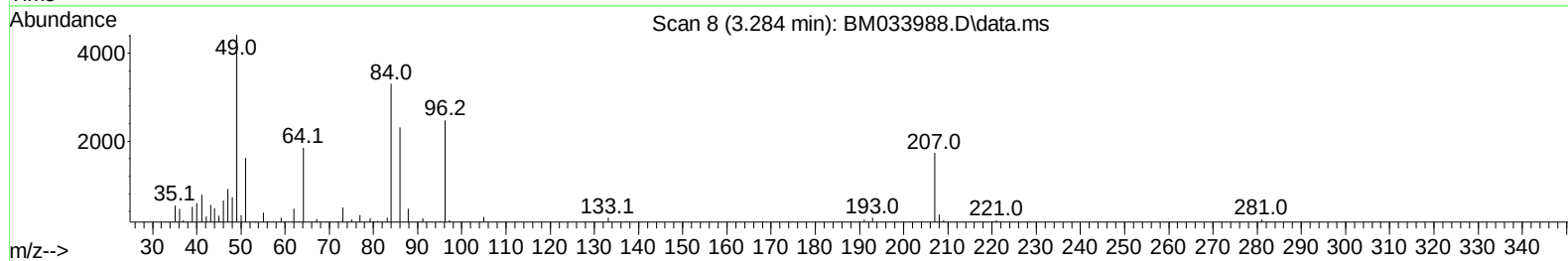
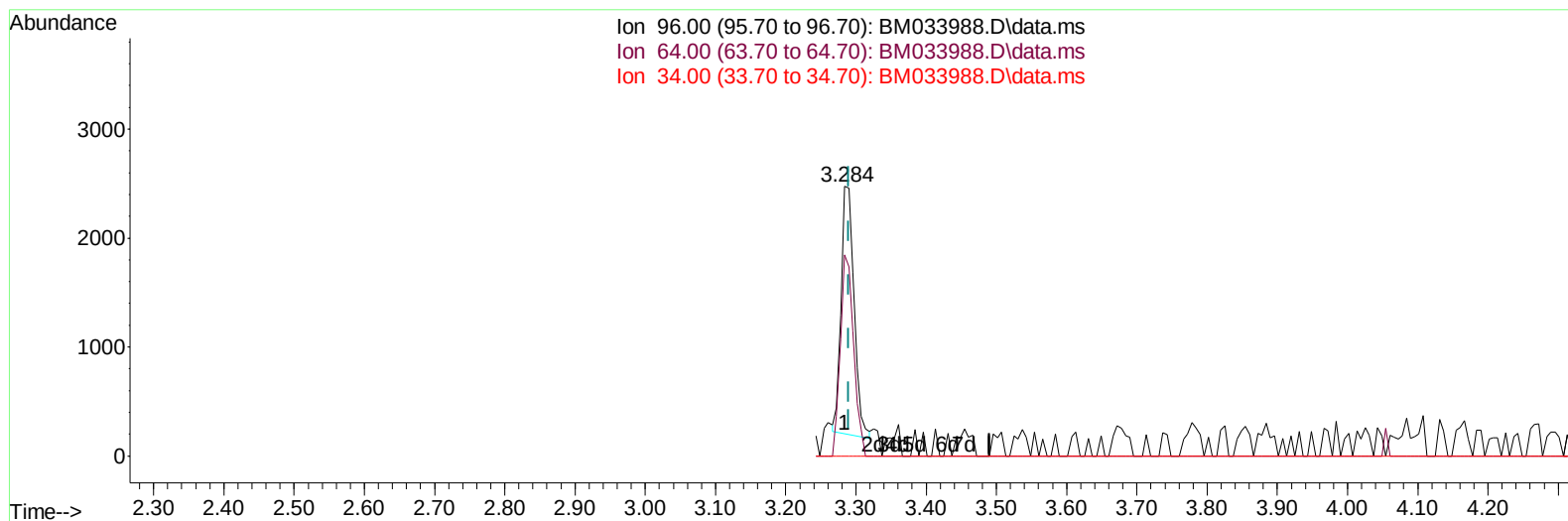
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020122\
 Data File : BM033988.D
 Acq On : 02 Feb 2022 23:46
 Operator : CG/JU
 Sample : N1355-09DL 5X
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0VF3DL

Manual IntegrationsAPPROVED

Quant Time: Feb 03 02:53:08 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Feb 01 02:16:28 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/03/2022
 Supervised By :Yogesh Patel 02/04/2022



TIC: BM033988.D\data.ms

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

3.284min (-0.006) 0.82 ng/uL

response 2914

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	74.61
34.00	0.00	0.00
0.00	0.00	0.00

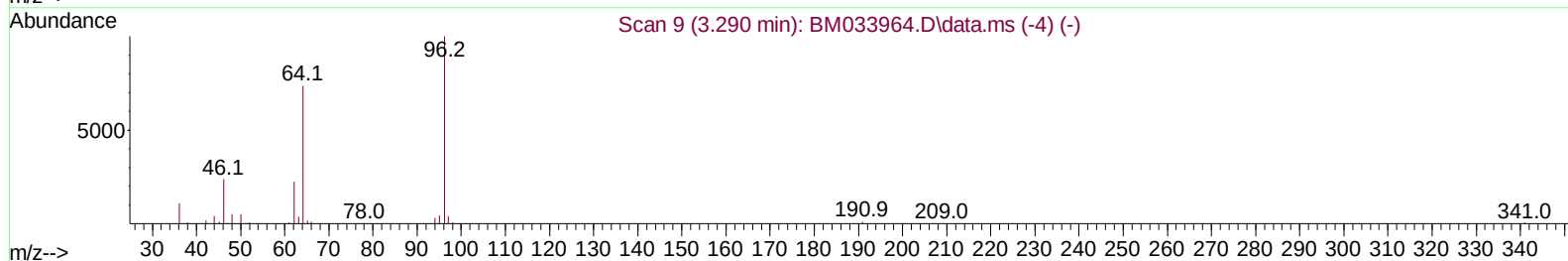
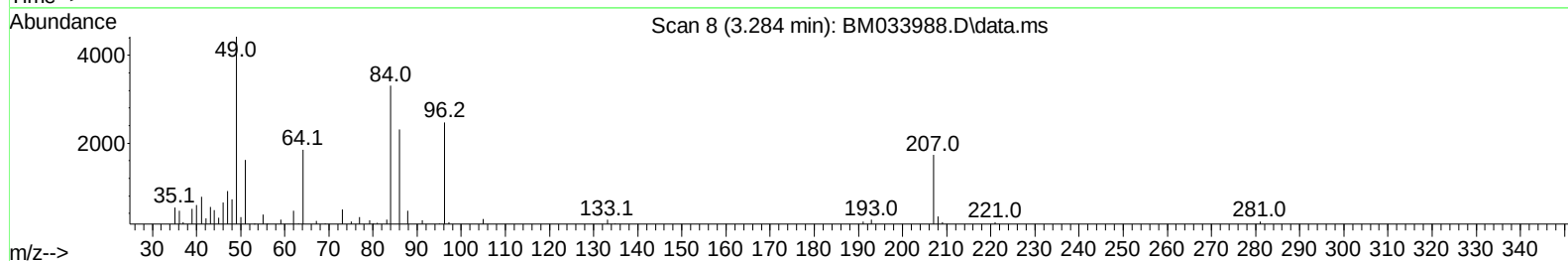
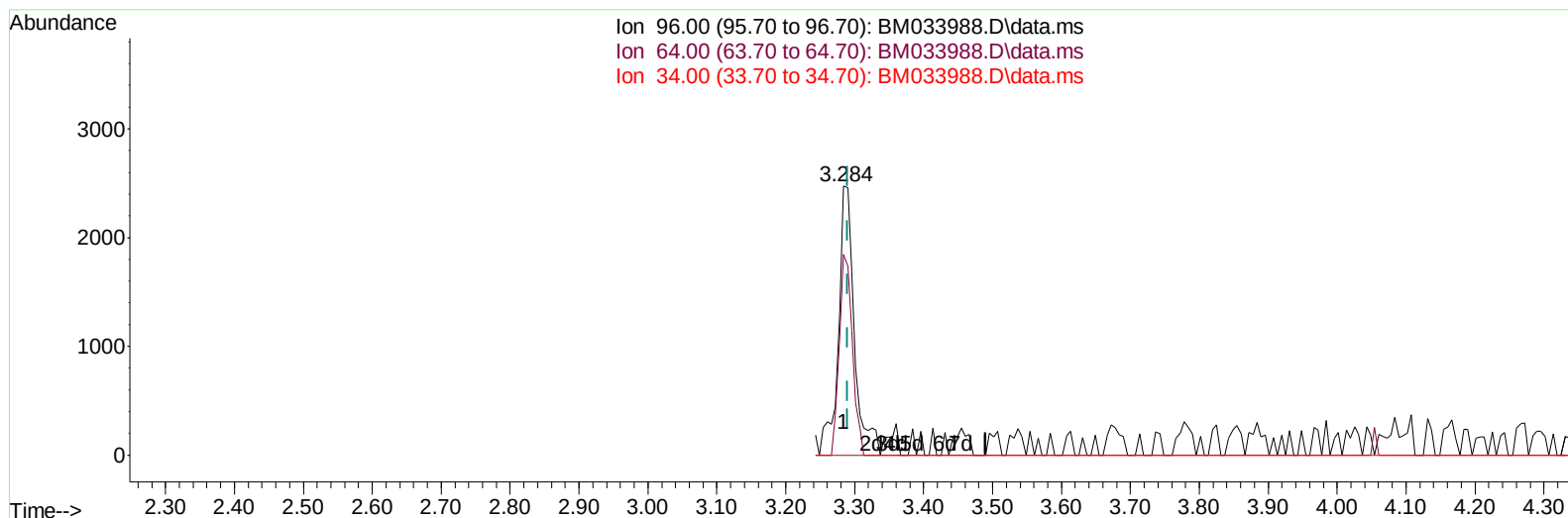
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TIC: BM033988.D\data.ms

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

3.284min (-0.006) 1.13 ng/uL m

response 3996

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	74.61
34.00	0.00	0.00
0.00	0.00	0.00

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 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :

BNA_M

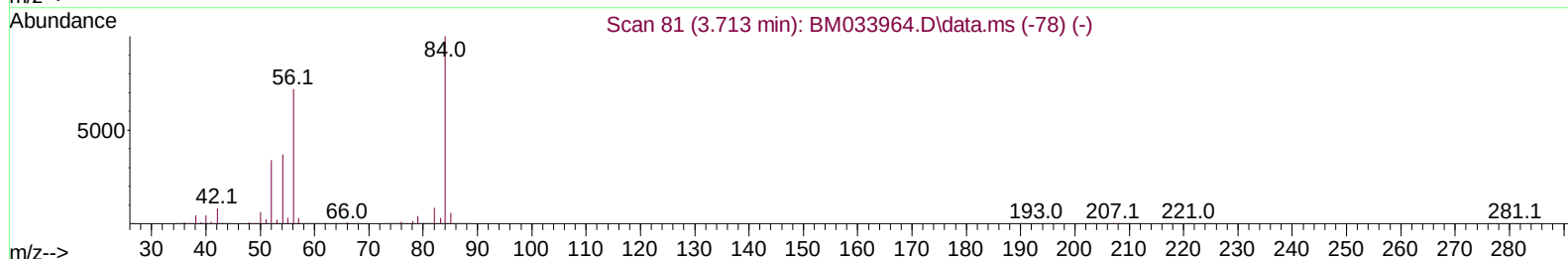
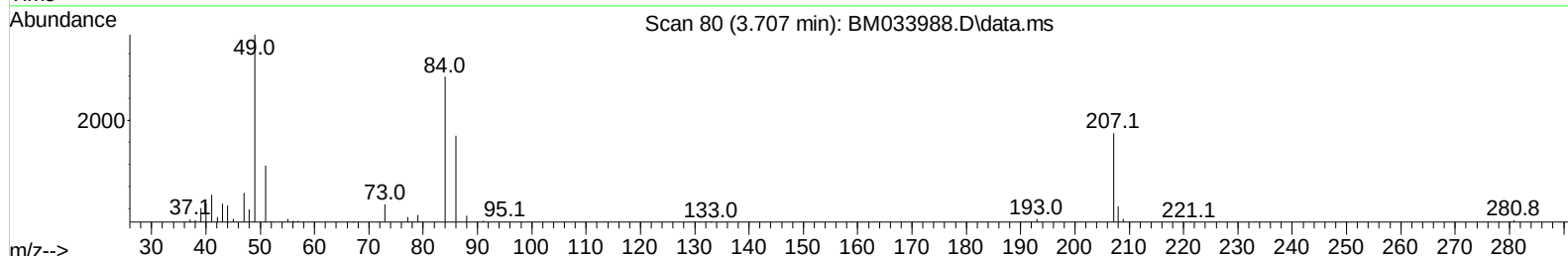
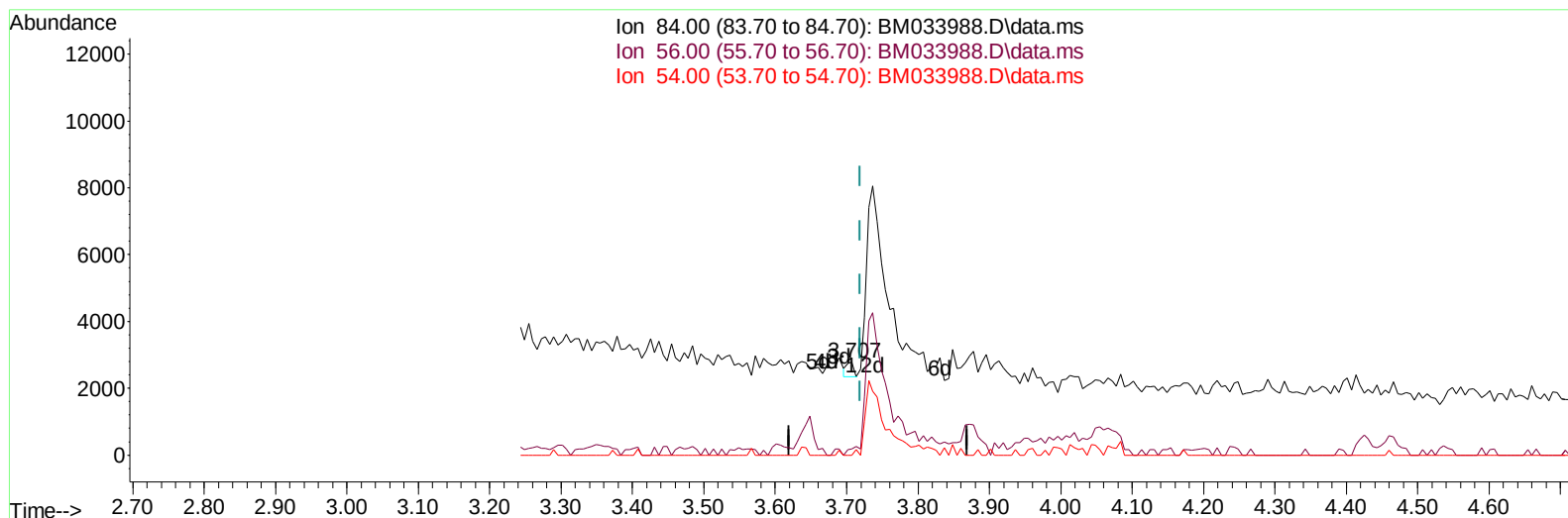
ClientSampleId :

C0VF3DL

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TIC: BM033988.D\data.ms

(4) Pyridine-d5 (S)

3.707min (-0.012) 0.03 ng/u1

response 301

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	80.70	6.25#
54.00	42.60	0.00#
0.00	0.00	0.00

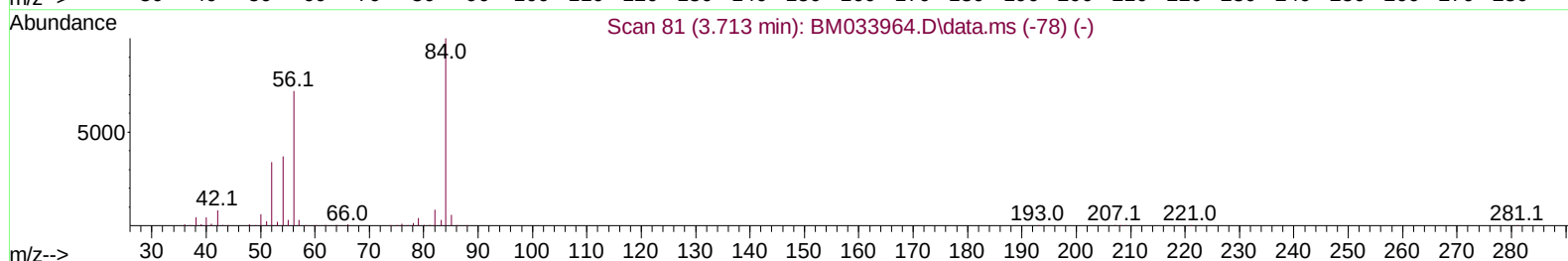
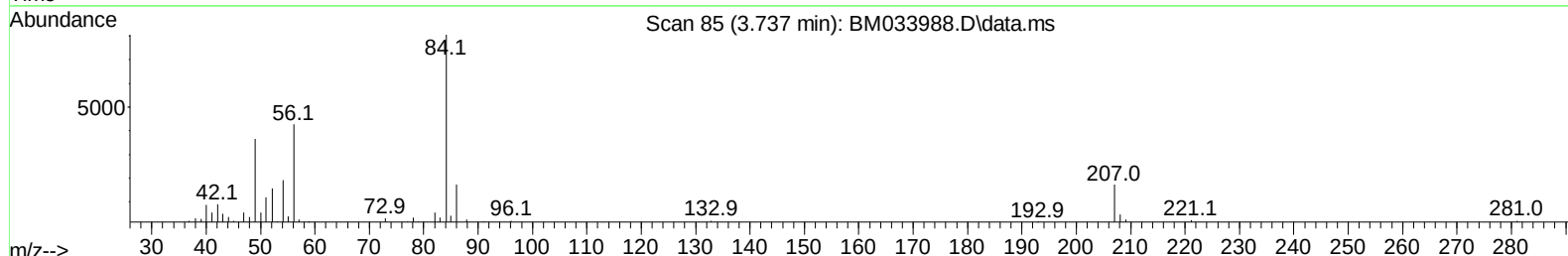
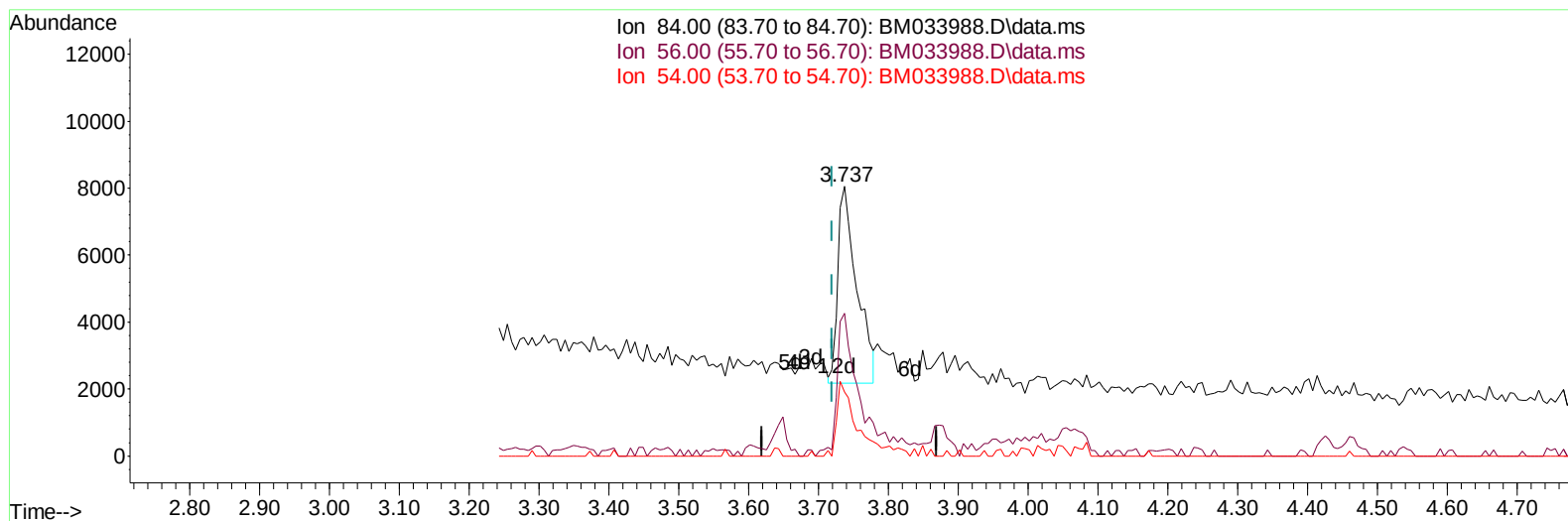
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 ALS Vial : 59 Sample Multiplier: 1

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TIC: BM033988.D\data.ms

(4) Pyridine-d5 (S)

3.737min (+ 0.018) 1.12 ng/ul m

response 11012

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	80.70	53.01#
54.00	42.60	23.81#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.925	152	139308	20.000	ng/ul	0.00
20) Naphthalene-d8	10.730	136	610576	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.560	164	364215	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.295	188	634032	20.000	ng/ul	-0.01
79) Chrysene-d12	21.459	240	492417	20.000	ng/ul	0.00
88) Perylene-d12	23.847	264	517486	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.284	96	3996m	1.131	ng/uL	0.00
4) Pyridine-d5	3.737	84	11012m	1.117	ng/ul	0.02
7) Phenol-d5	7.078	99	68055	5.831	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.242	67	42064	5.912	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.448	132	53971	5.772	ng/ul	-0.01
15) 4-Methylphenol-d8	8.631	113	53999	5.839	ng/ul	-0.01
21) Nitrobenzene-d5	9.084	128	27192	5.550	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.813	143	28645	5.414	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.348	165	53375	5.575	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.866	131	50878	3.848	ng/ul	0.00
46) Dimethylphthalate-d6	13.966	166	158073	5.808	ng/ul	-0.01
49) Acenaphthylene-d8	14.254	160	186125	5.583	ng/ul	0.00
54) 4-Nitrophenol-d4	14.760	143	13366	2.874	ng/ul	0.00
60) Fluorene-d10	15.548	176	129199	5.685	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.660	200	13072	3.284	ng/ul	-0.01
73) Anthracene-d10	17.395	188	171119	5.636	ng/ul	0.00
81) Pyrene-d10	19.677	212	177371	6.101	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.688	264	153207	5.637	ng/ul	-0.02

Target Compounds Qvalue

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

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