

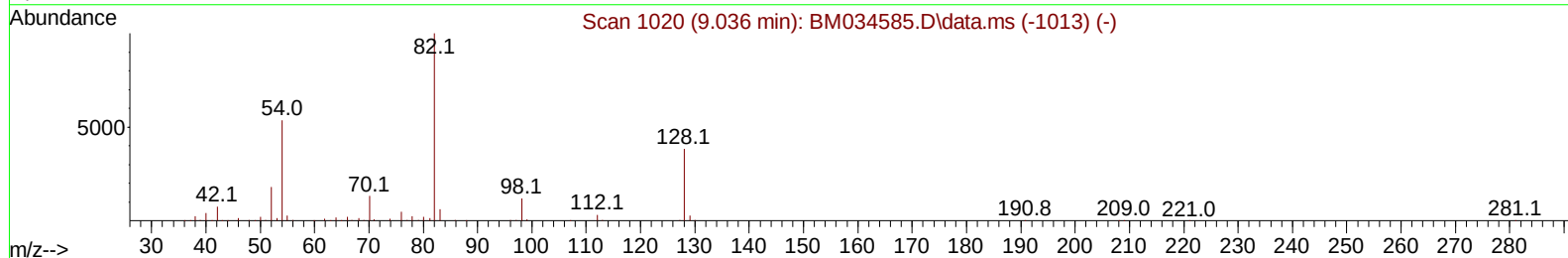
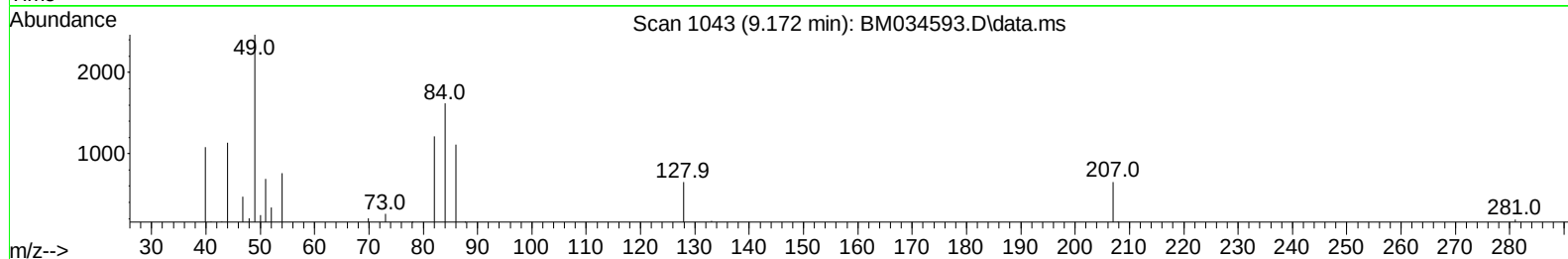
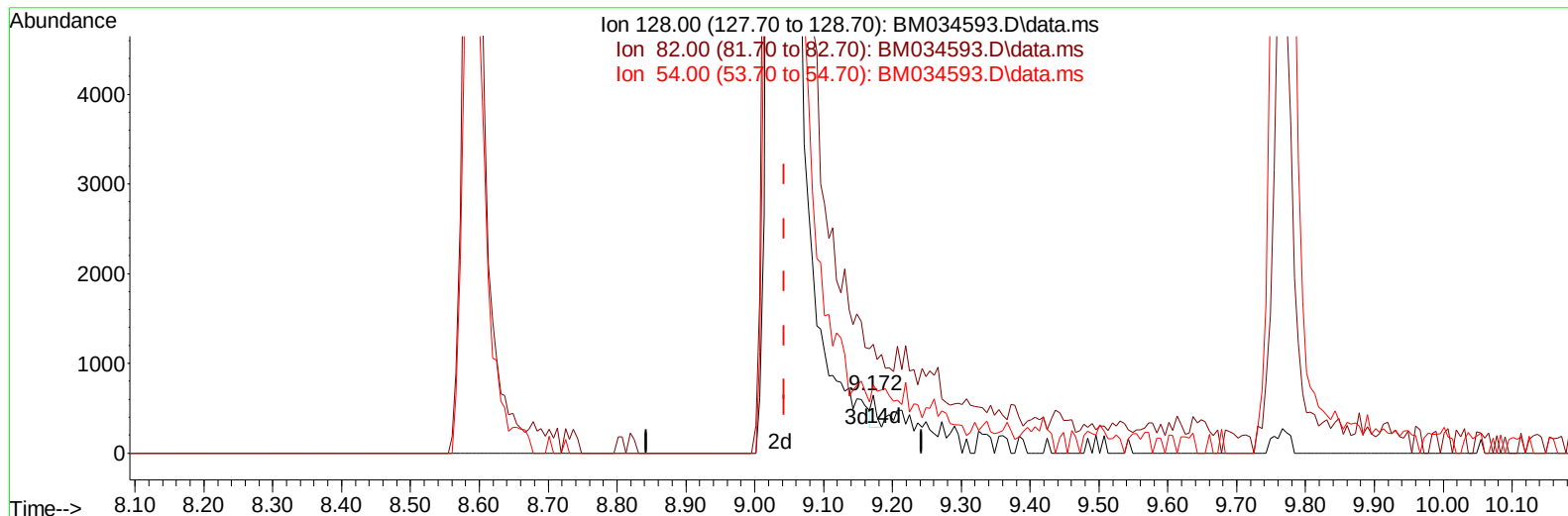
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040922\
Data File : BM034593.D
Acq On : 09 Apr 2022 11:30
Operator : CG/JU
Sample : N2319-03
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
YB2H0

Manual IntegrationsAPPROVED

Quant Time: Apr 11 00:59:07 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Apr 08 14:30:08 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/11/2022
Supervised By :Yogesh Patel 04/12/2022



TIC: BM034593.D\data.ms

(21) Nitrobenzene-d5 (S)

9.172min (+ 0.130) 0.06 ng/ul

response 167

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	228.00	187.64
54.00	100.50	117.00
0.00	0.00	0.00

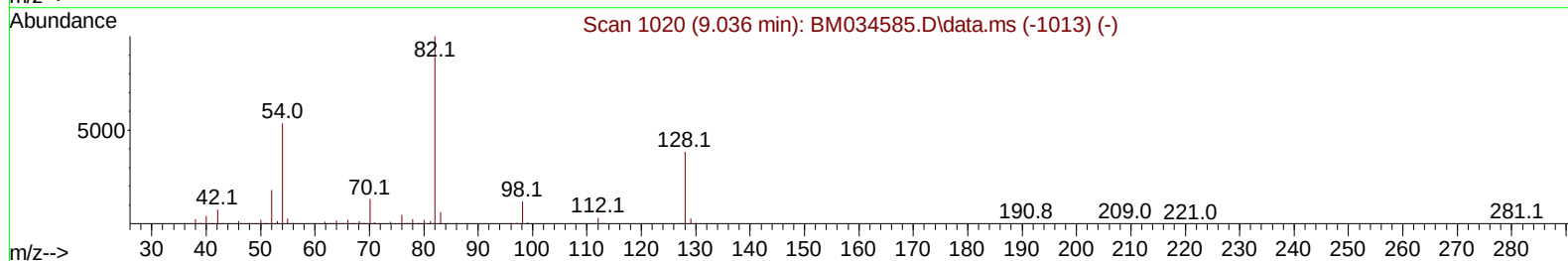
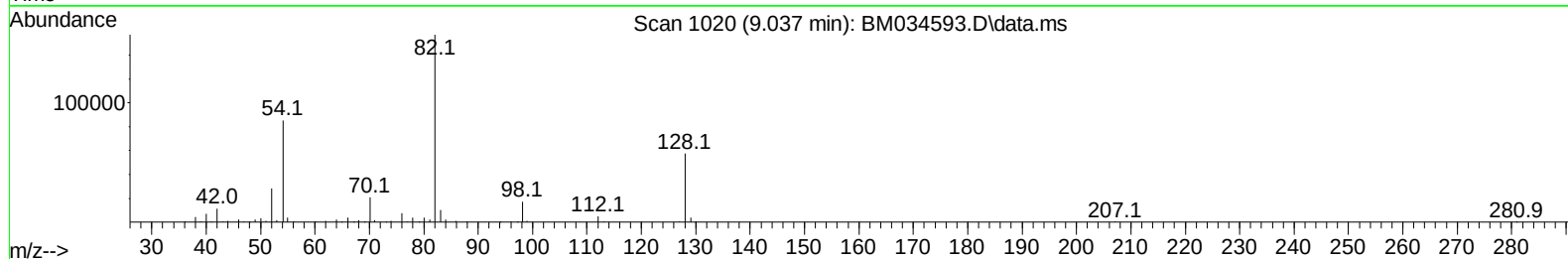
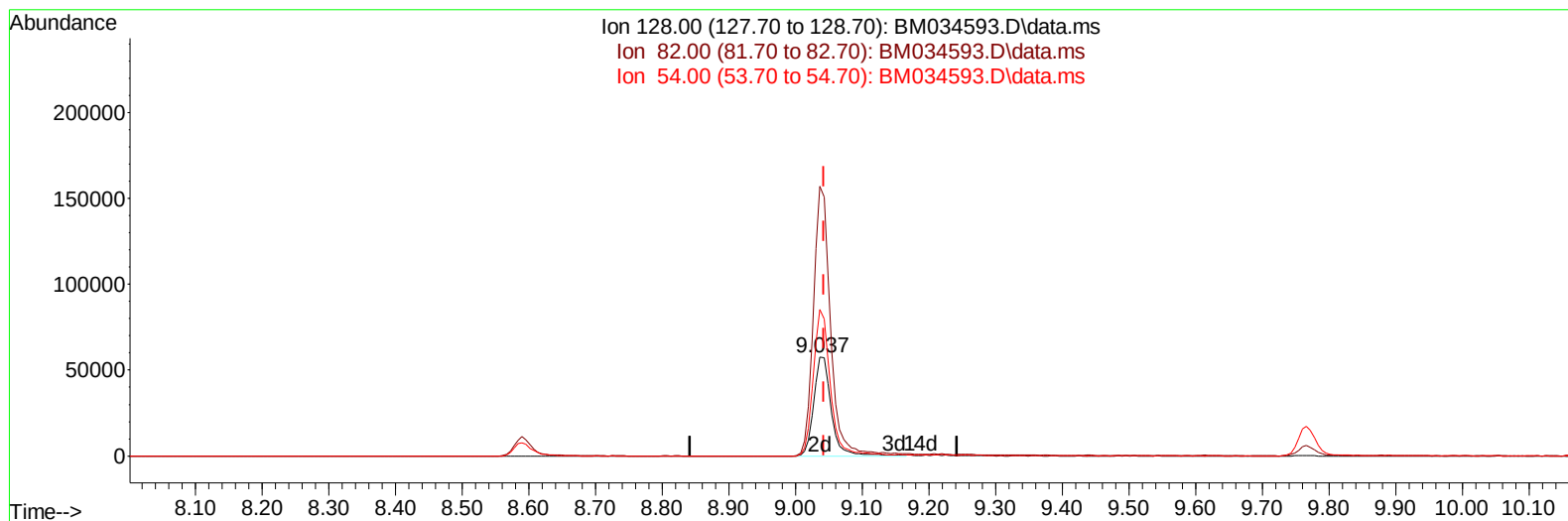
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(21) Nitrobenzene-d5 (S)

9.037min (-0.006) 35.60 ng/ul m

response 104681

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	228.00	272.82
54.00	100.50	148.29#
0.00	0.00	0.00

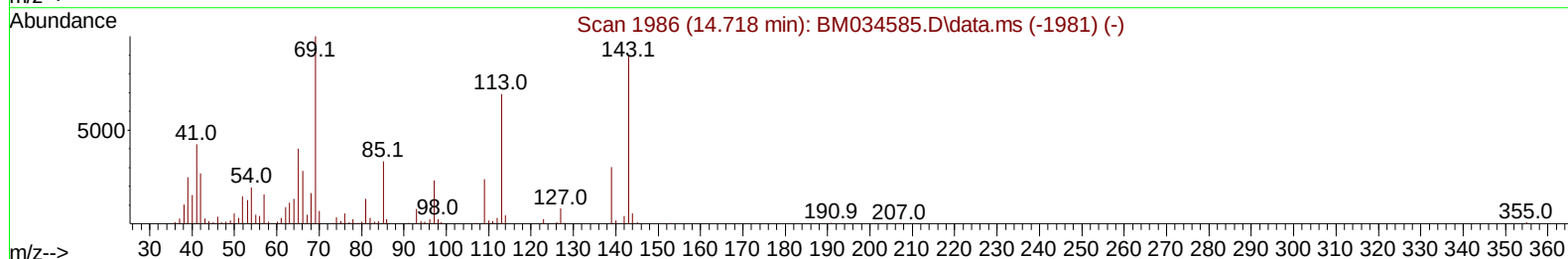
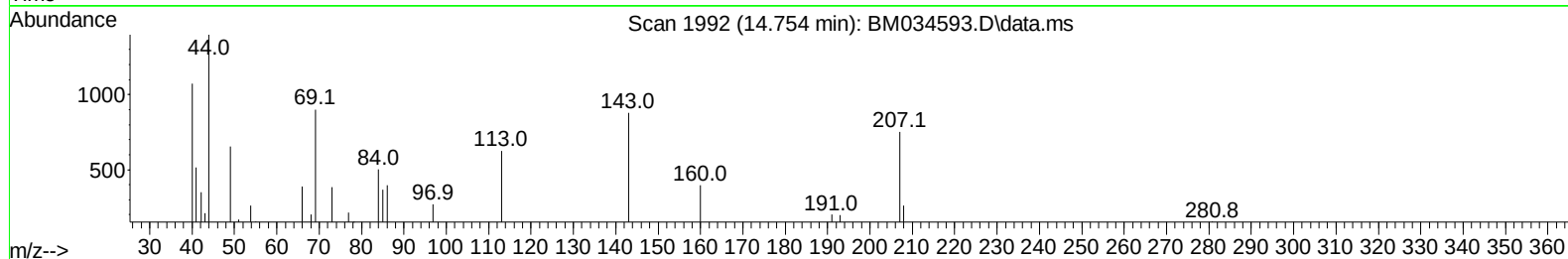
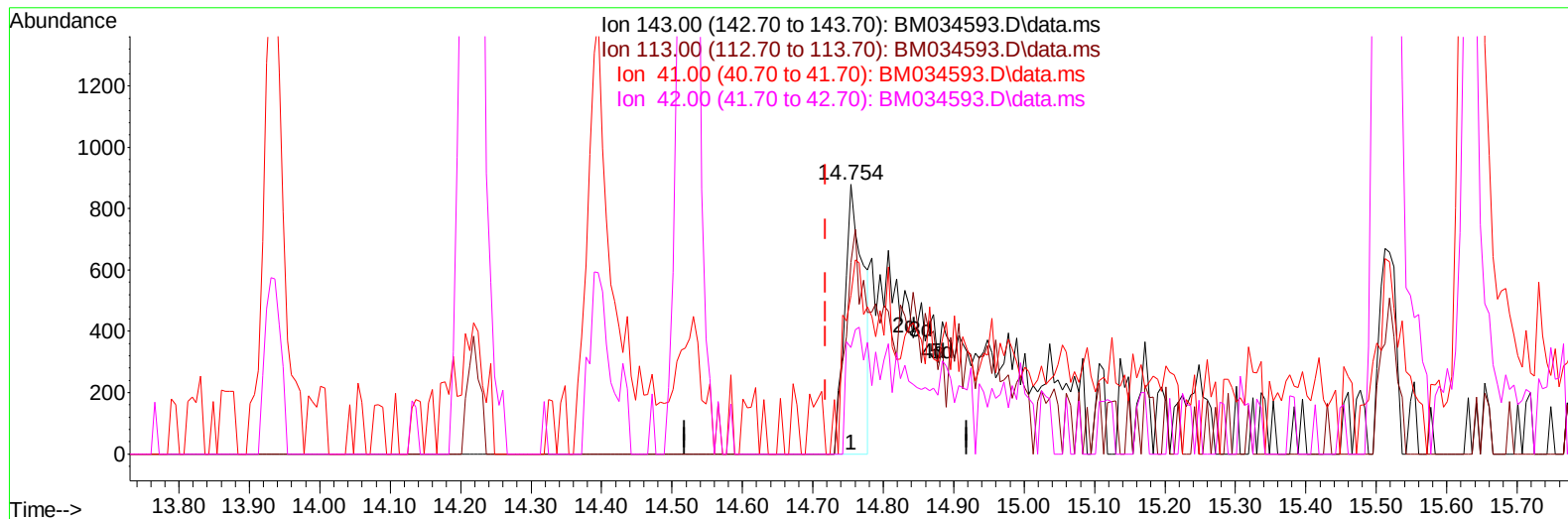
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Instrument :
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TIC: BM034593.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.754min (+ 0.036) 0.65 ng/ul

response 1615

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	71.10
41.00	31.90	58.82#
42.00	19.10	39.59#

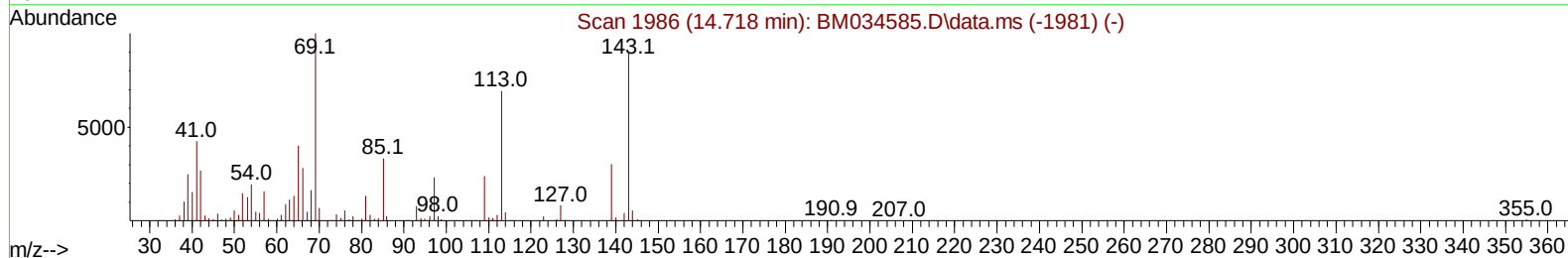
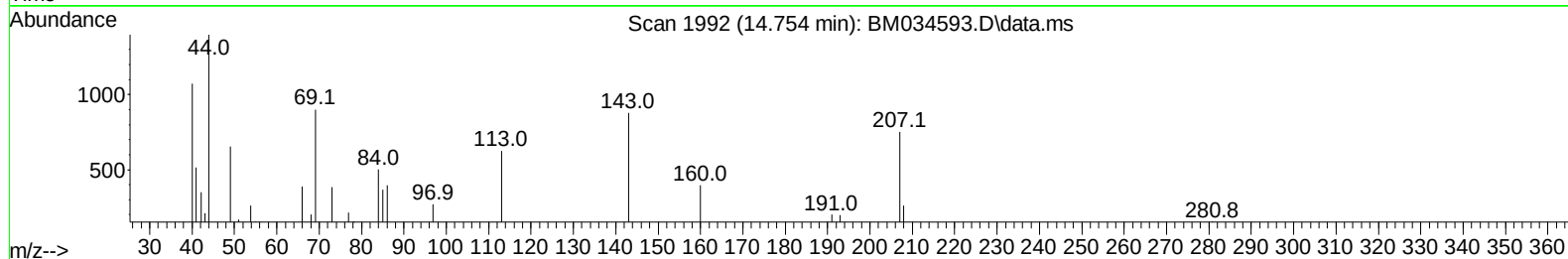
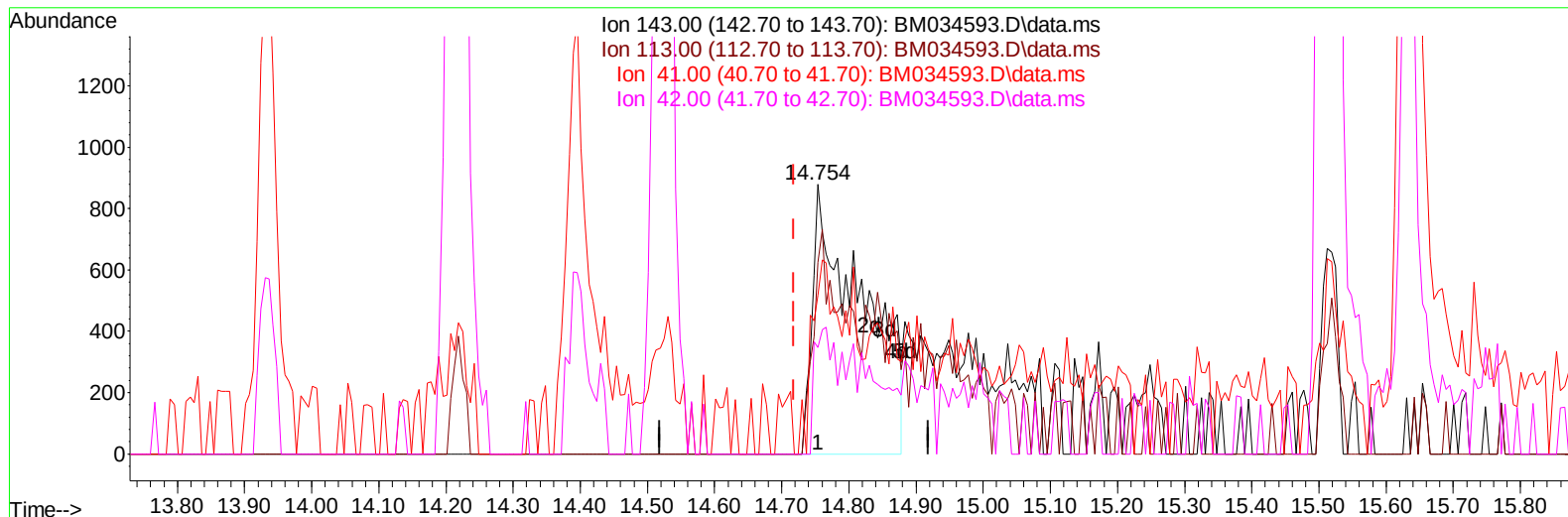
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Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
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TIC: BM034593.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.754min (+ 0.036) 1.82 ng/ul m

response 4498

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	71.10
41.00	31.90	58.82#
42.00	19.10	39.59#

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 ALS Vial : 40 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.878	152	100627	20.000	ng/ul	0.00
20) Naphthalene-d8	10.684	136	400119	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.525	164	223848	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.266	188	419296	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.442	240	407698	20.000	ng/ul	# 0.00
88) Perylene-d12	23.824	264	492131	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.249	96	12740	4.515	ng/uL	0.00
4) Pyridine-d5	3.684	84	52173	7.661	ng/ul	0.00
7) Phenol-d5	7.043	99	45001	5.373	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.202	67	191787	34.364	ng/ul	0.00
11) 2-Chlorophenol-d4	7.407	132	167853	25.925	ng/ul	0.00
15) 4-Methylphenol-d8	8.590	113	93034	14.220	ng/ul	0.00
21) Nitrobenzene-d5	9.037	128	104681m	35.598	ng/ul	0.00
24) 2-Nitrophenol-d4	9.766	143	101483	33.147	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.307	165	154273	26.760	ng/ul	0.00
31) 4-Chloroaniline-d4	10.825	131	205220	25.803	ng/ul	0.00
46) Dimethylphthalate-d6	13.936	166	578752	35.781	ng/ul	0.00
49) Acenaphthylene-d8	14.219	160	721421	34.702	ng/ul	0.00
54) 4-Nitrophenol-d4	14.754	143	4498m	1.824	ng/ul	0.04
60) Fluorene-d10	15.513	176	532308	38.083	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.636	200	79610	31.101	ng/ul	0.00
73) Anthracene-d10	17.366	188	866612	43.351	ng/ul	0.00
81) Pyrene-d10	19.654	212	1061852	47.719	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.671	264	1052573	42.519	ng/ul	0.00

Target Compounds Qvalue

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

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