

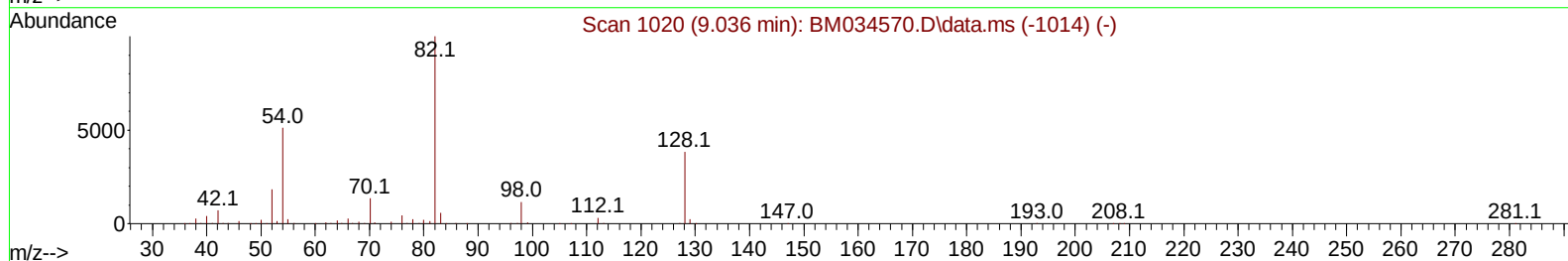
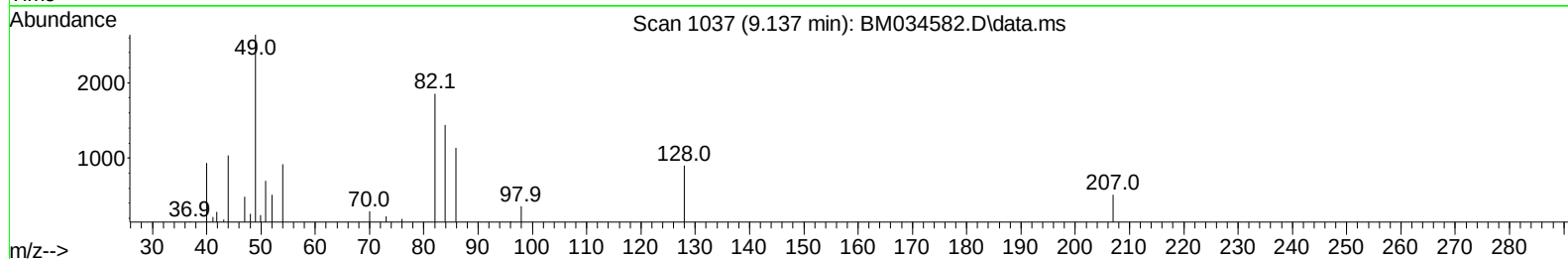
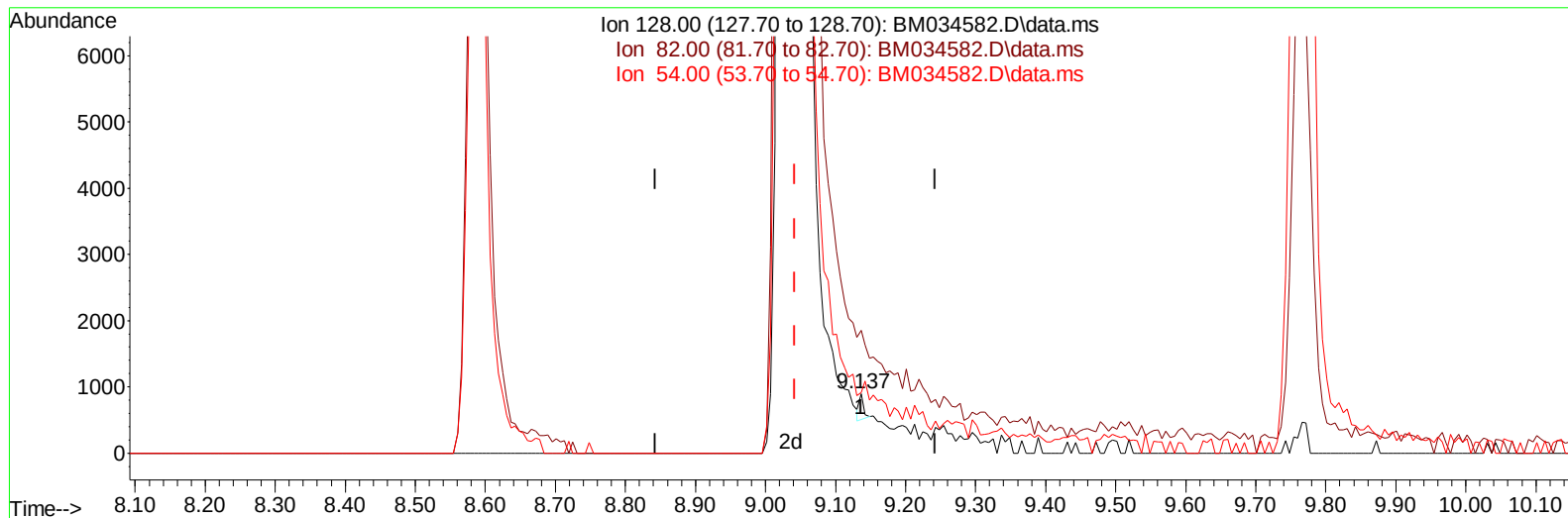
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040922\
Data File : BM034582.D
Acq On : 09 Apr 2022 04:14
Operator : CG/JU
Sample : N2319-04
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
YB2H1

Manual IntegrationsAPPROVED

Quant Time: Apr 09 04:48:01 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: BM034582.D\data.ms

(21) Nitrobenzene-d5 (S)

9.137min (+ 0.094) 0.05 ng/u1

response 166

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	228.00	206.00
54.00	100.50	101.56
0.00	0.00	0.00

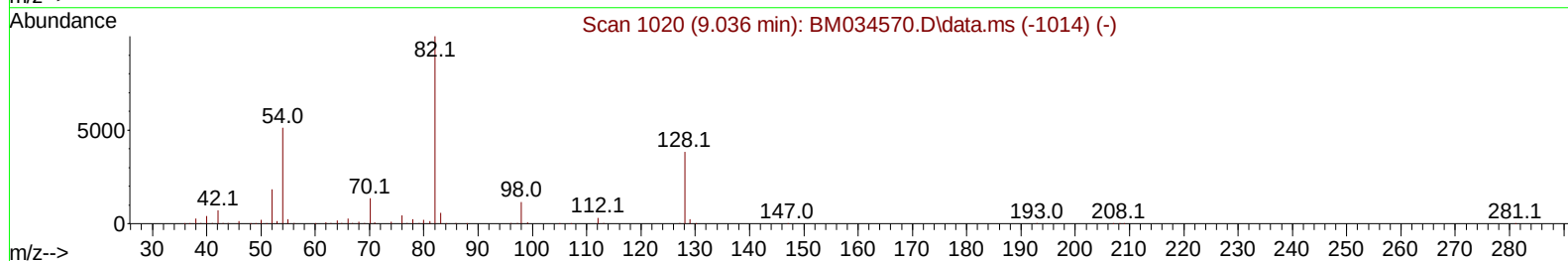
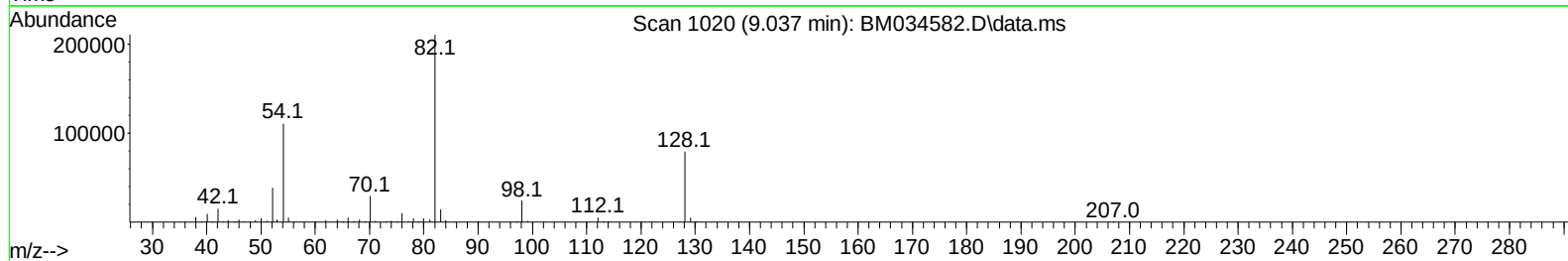
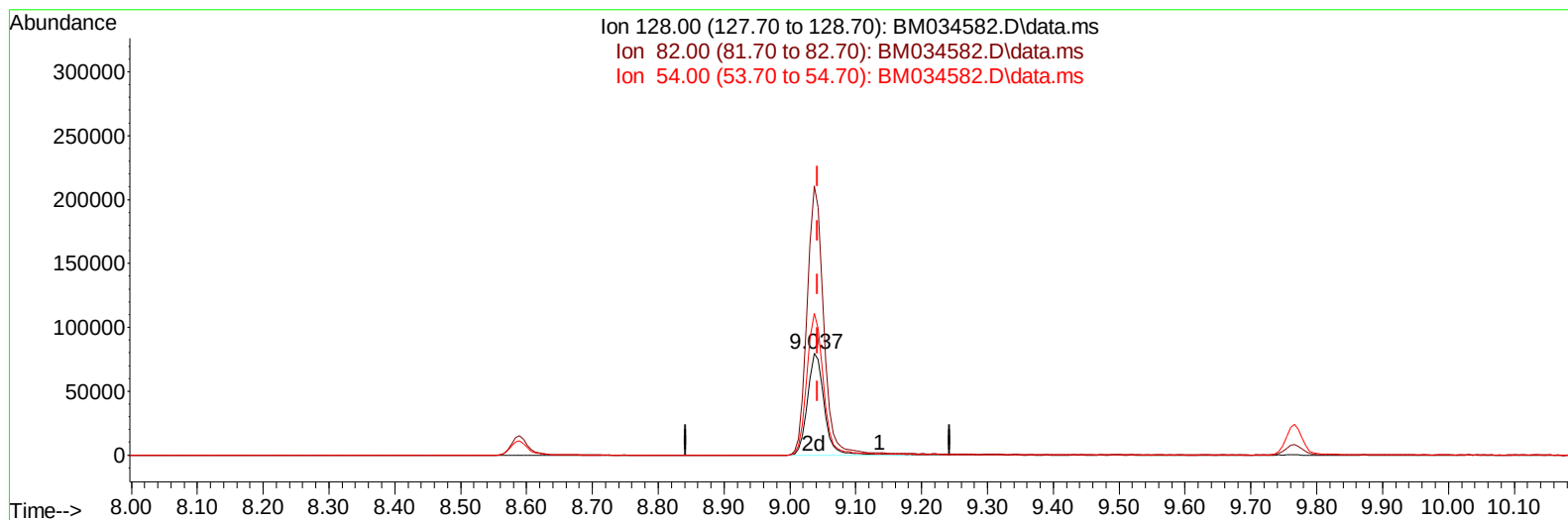
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TIC: BM034582.D\data.ms

(21) Nitrobenzene-d5 (S)

9.037min (-0.006) 38.58 ng/ul m

response 139632

Ion	Exp%	Act%
128.00	100.00	100.00
82.00	228.00	264.28
54.00	100.50	139.22#
0.00	0.00	0.00

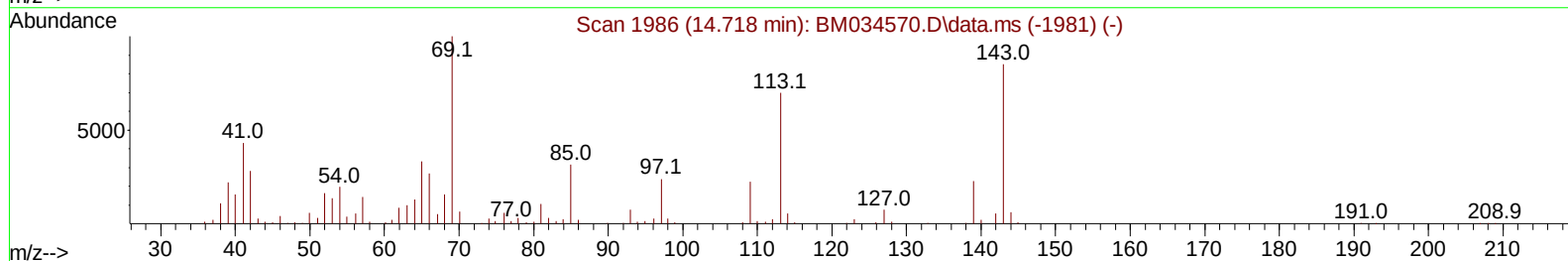
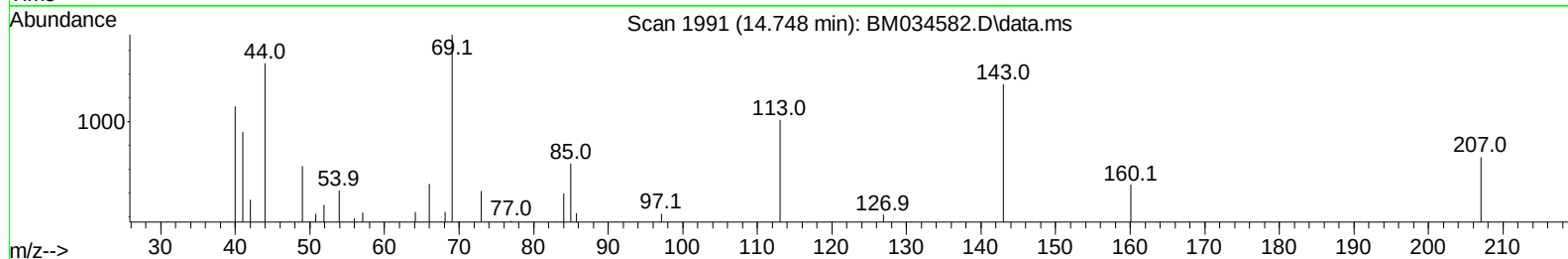
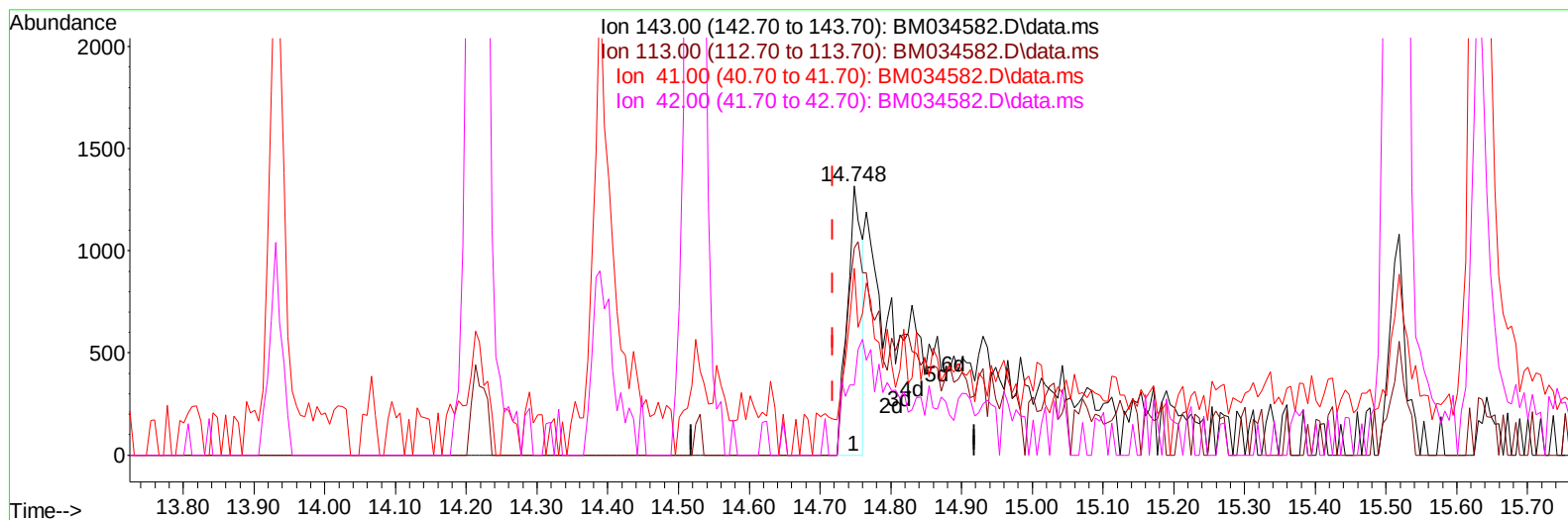
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 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
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TIC: BM034582.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.748min (+ 0.030) 0.63 ng/ul

response 1895

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	76.99
41.00	31.90	69.48#
42.00	19.10	26.04#

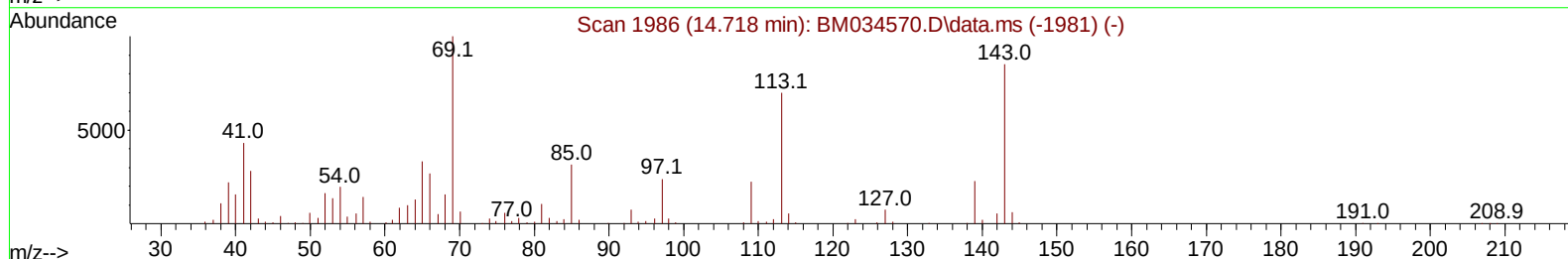
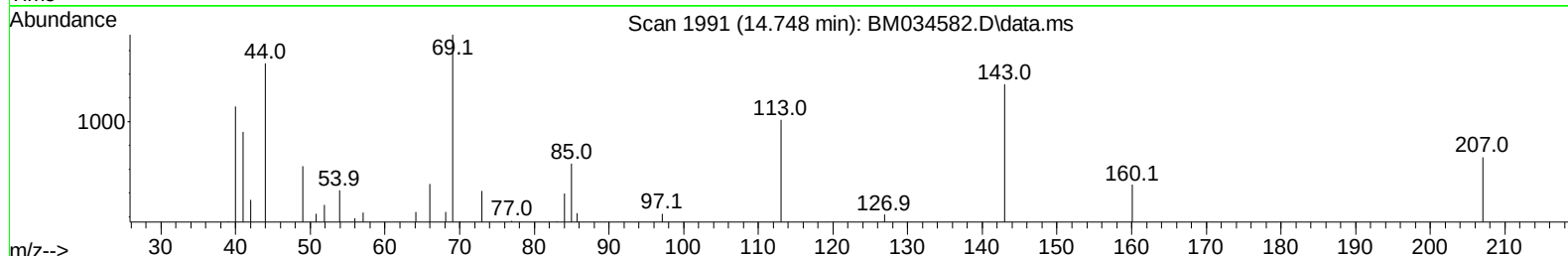
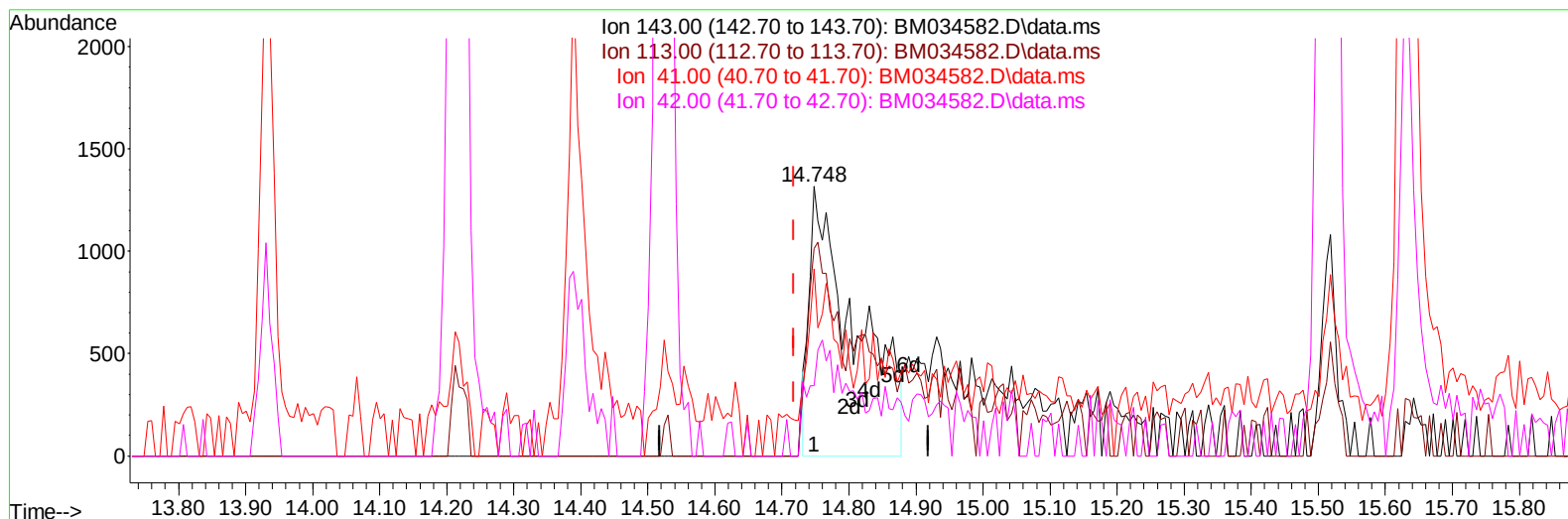
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 Sample : N2319-04
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
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TIC: BM034582.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.748min (+ 0.030) 2.10 ng/ul m

response 6269

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	76.99
41.00	31.90	69.48#
42.00	19.10	26.04#

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 Misc :
 ALS Vial : 28 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	124333	20.000	ng/ul	0.00
20) Naphthalene-d8	10.684	136	492397	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.524	164	271586	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.265	188	515656	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.442	240	476370	20.000	ng/ul	# 0.00
88) Perylene-d12	23.818	264	564671	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.255	96	16237	4.657	ng/uL	0.00
4) Pyridine-d5	3.684	84	64349	7.647	ng/ul	0.00
7) Phenol-d5	7.037	99	61589	5.952	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.201	67	254854	36.958	ng/ul	0.00
11) 2-Chlorophenol-d4	7.407	132	225532	28.192	ng/ul	0.00
15) 4-Methylphenol-d8	8.590	113	124475	15.398	ng/ul	0.00
21) Nitrobenzene-d5	9.037	128	139632m	38.585	ng/ul	0.00
24) 2-Nitrophenol-d4	9.766	143	136934	36.344	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.307	165	218263	30.765	ng/ul	0.00
31) 4-Chloroaniline-d4	10.825	131	270759	27.663	ng/ul	0.00
46) Dimethylphthalate-d6	13.936	166	792078	40.362	ng/ul	0.00
49) Acenaphthylene-d8	14.219	160	943293	37.398	ng/ul	0.00
54) 4-Nitrophenol-d4	14.748	143	6269m	2.095	ng/ul	0.03
60) Fluorene-d10	15.513	176	681767	40.202	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.630	200	102181	32.459	ng/ul	0.00
73) Anthracene-d10	17.365	188	1060700	43.144	ng/ul	0.00
81) Pyrene-d10	19.654	212	1274838	49.032	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.665	264	1158267	40.778	ng/ul	0.00

Target Compounds Qvalue

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

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