

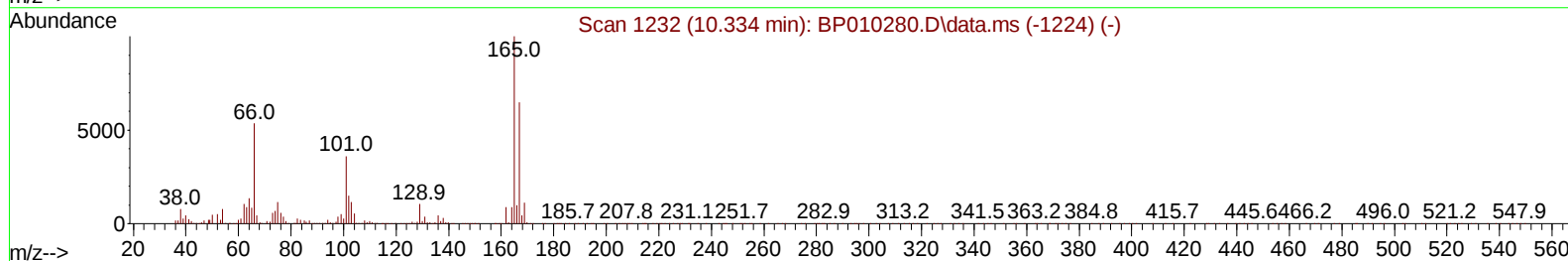
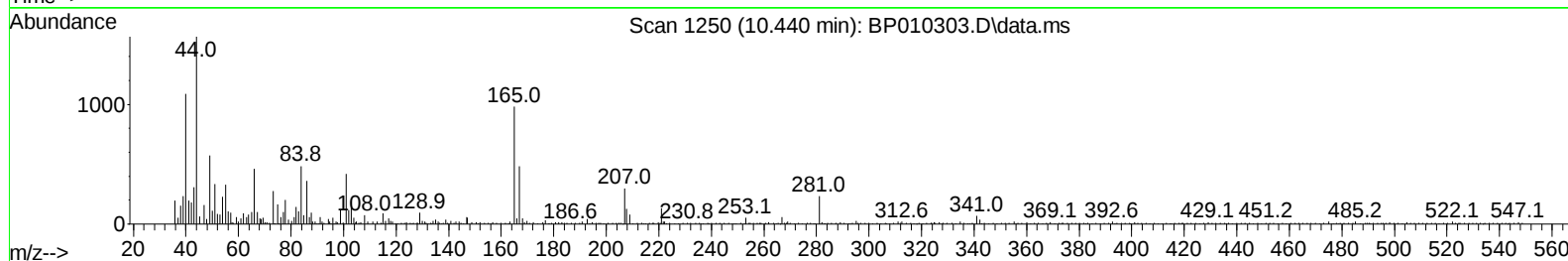
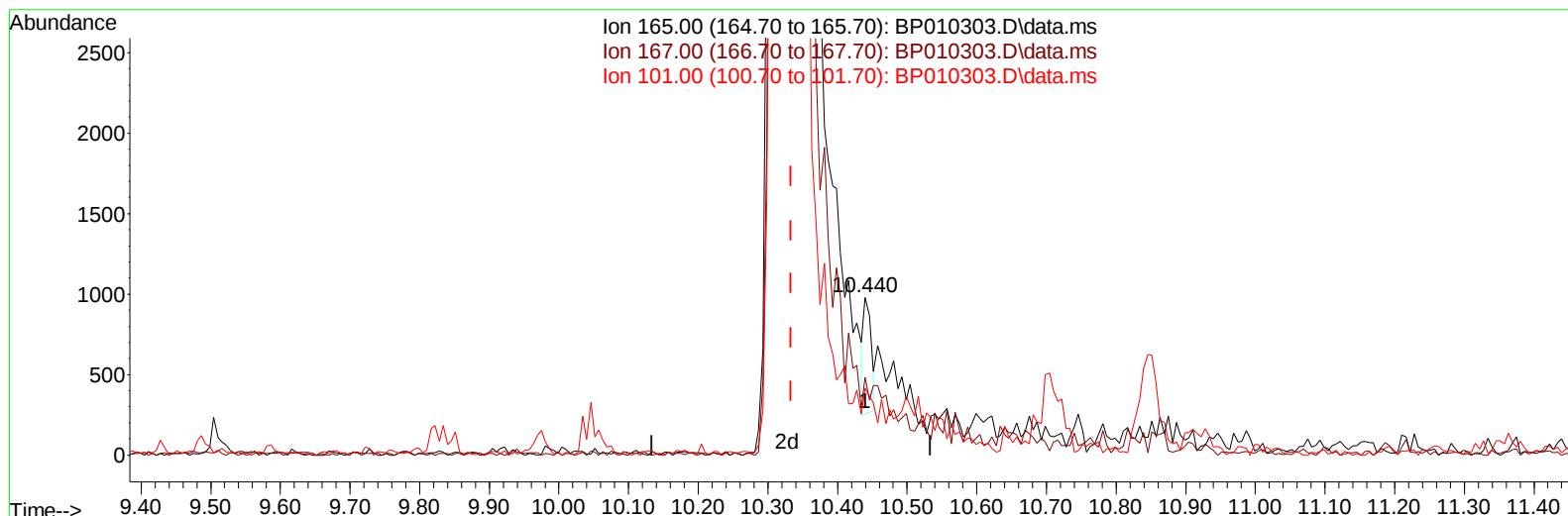
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
Data File : BP010303.D
Acq On : 12 May 2022 01:35
Operator : CG/JU
Sample : N2707-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EXL83

Manual IntegrationsAPPROVED

Quant Time: May 12 02:59:31 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: BP010303.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.440min (+ 0.106) 0.04 ng/u1

response 558

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	49.39
101.00	49.80	42.55
0.00	0.00	0.00

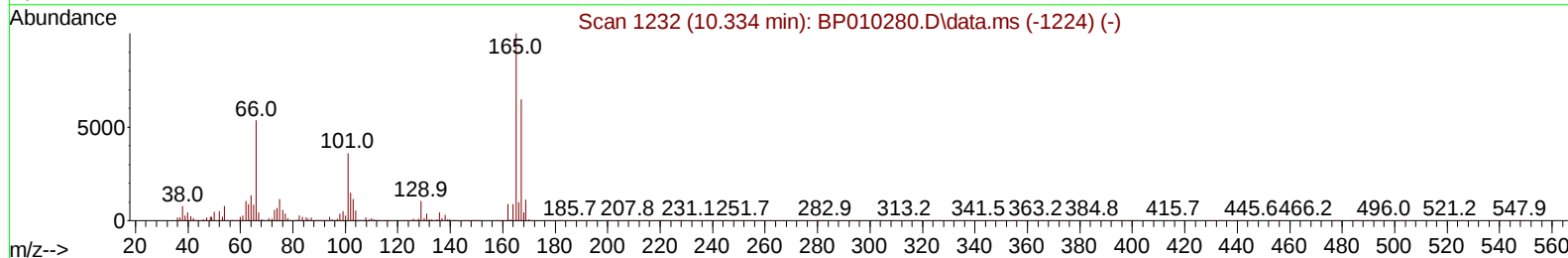
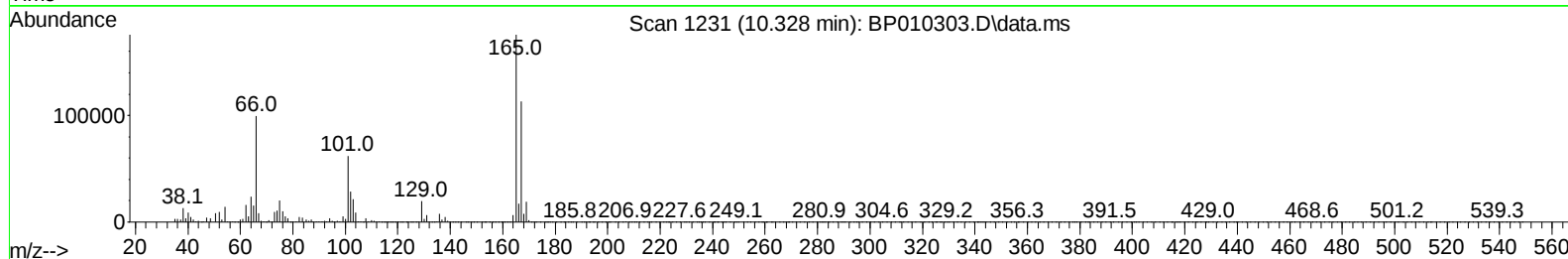
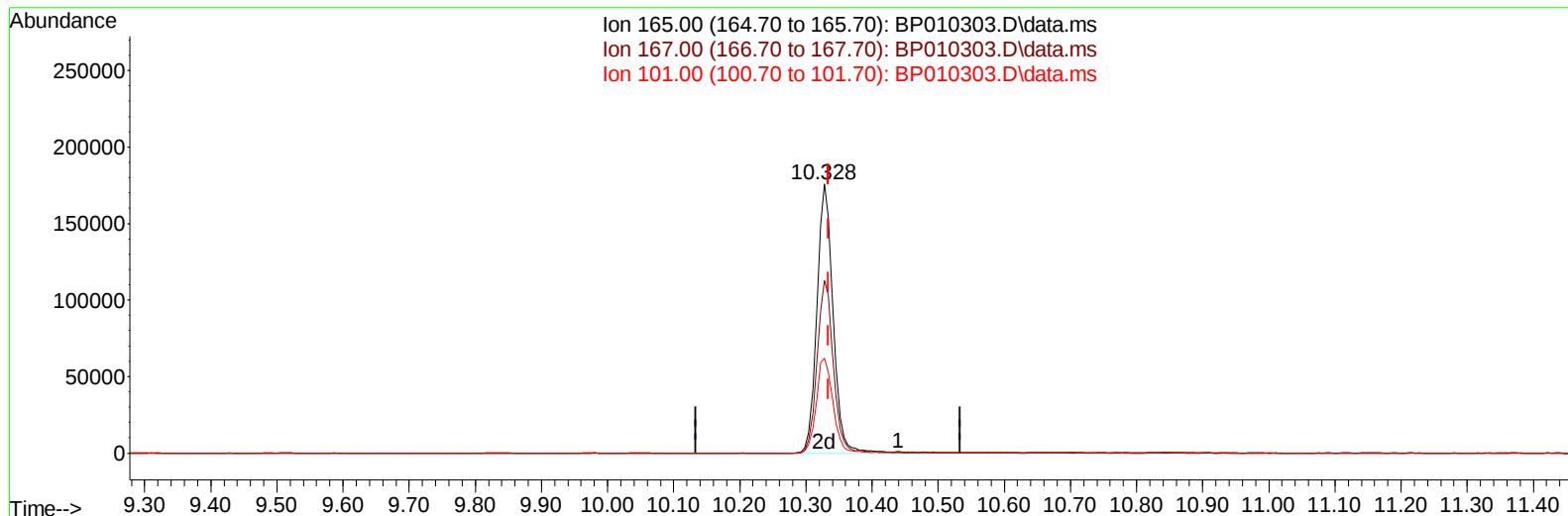
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TIC: BP010303.D\data.ms

(28) 2,4-Dichlorophenol-d3 (S)

10.328min (-0.006) 23.47 ng/ul m

response 302236

Ion	Exp%	Act%
165.00	100.00	100.00
167.00	57.70	64.29
101.00	49.80	35.31#
0.00	0.00	0.00

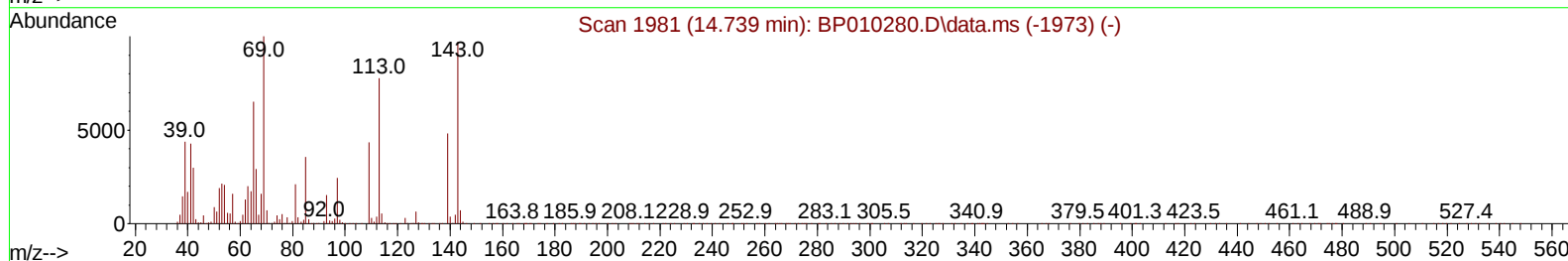
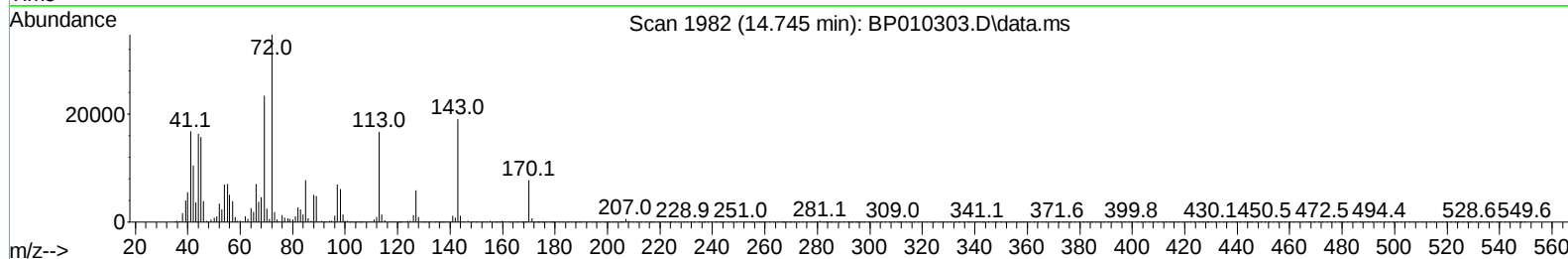
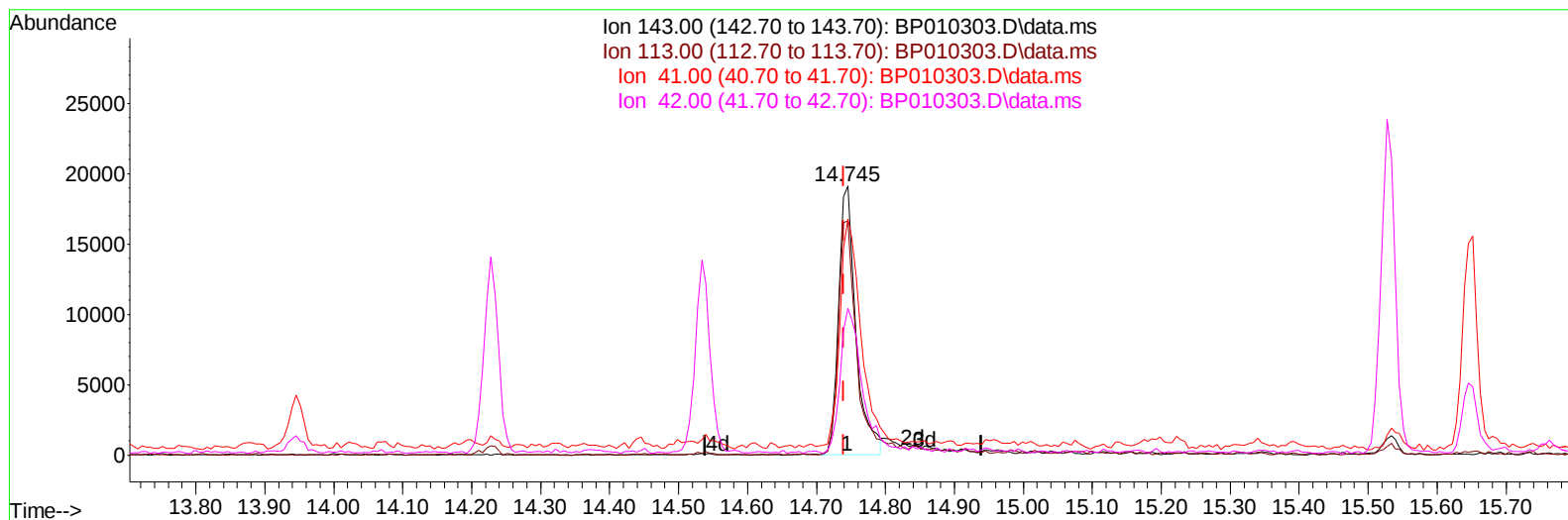
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Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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TIC: BP010303.D\data.ms

(54) 4-Nitrophenol-d4 (S)**14.745min (+ 0.006) 3.98 ng/ul****response 33241**

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	87.23#
41.00	23.20	87.82#
42.00	18.00	54.53#

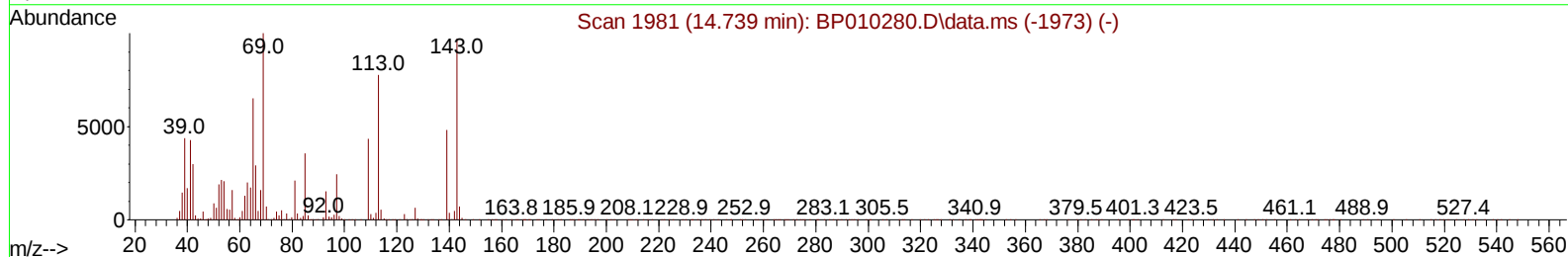
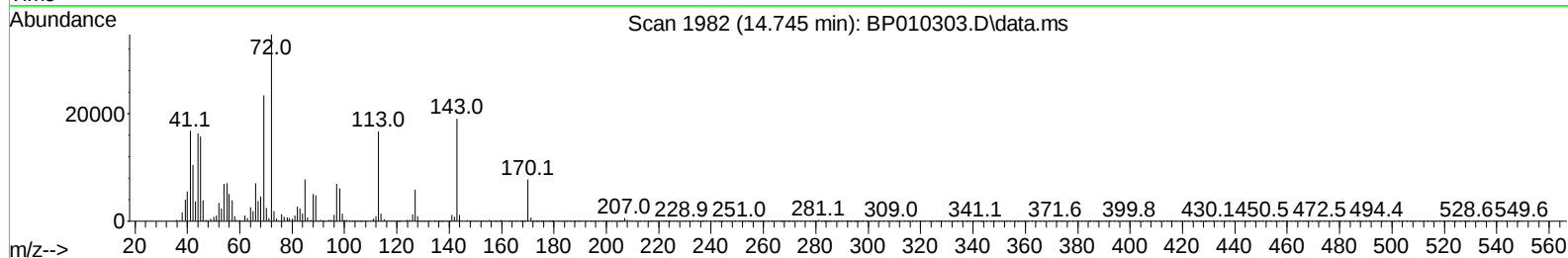
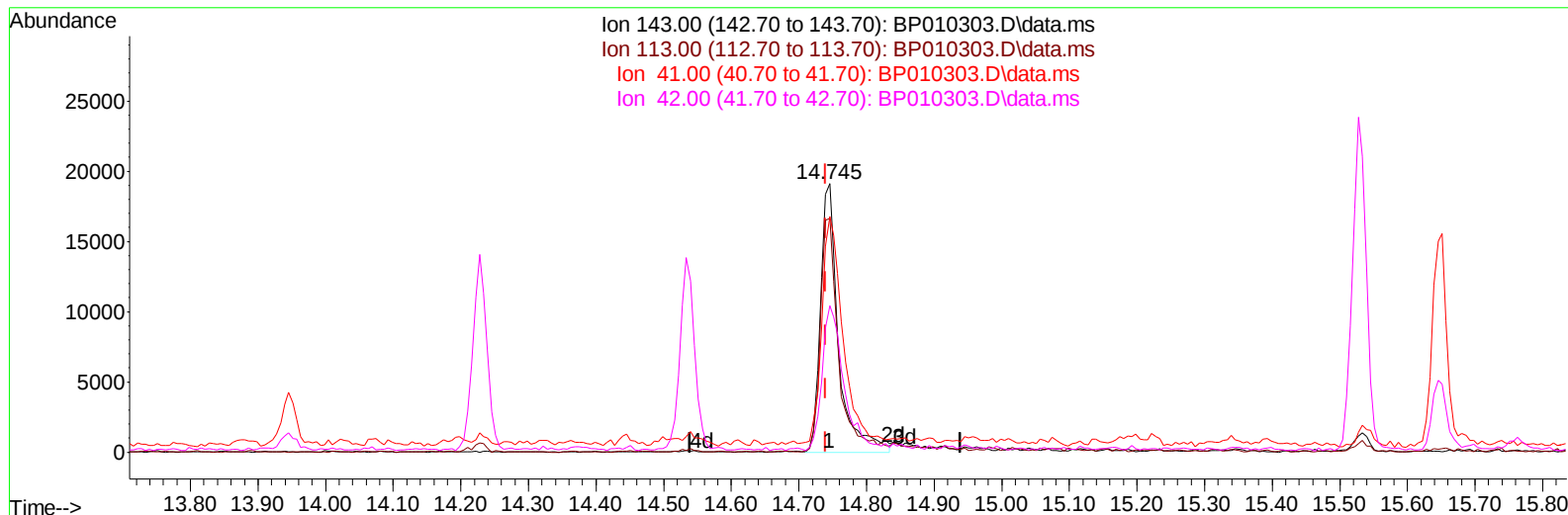
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Operator : CG/JU
Sample : N2707-01
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EXL83

Manual IntegrationsAPPROVED

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TIC: BP010303.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.745min (+ 0.006) 4.26 ng/ul m

response 35561

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	87.23#
41.00	23.20	87.82#
42.00	18.00	54.53#

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 Misc :
 ALS Vial : 26 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.905	152	178540	20.000	ng/ul	0.00
20) Naphthalene-d8	10.704	136	774002	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.534	164	546319	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.286	188	1195785	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.369	240	1168147	20.000	ng/ul	# 0.00
88) Perylene-d12	23.798	264	1065326	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.352	96	19900	4.735	ng/uL	0.00
4) Pyridine-d5	3.781	84	32363	2.767	ng/ul	0.01
7) Phenol-d5	7.069	99	77058	4.647	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	286858	27.963	ng/ul	0.00
11) 2-Chlorophenol-d4	7.440	132	251551	20.598	ng/ul	0.00
15) 4-Methylphenol-d8	8.610	113	166030	12.178	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128	179924	28.992	ng/ul	0.00
24) 2-Nitrophenol-d4	9.787	143	185041	26.151	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165	302236m	23.467	ng/ul	0.00
31) 4-Chloroaniline-d4	10.840	131	351734	18.499	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	1251577	28.756	ng/ul	0.00
49) Acenaphthylene-d8	14.228	160	1395314	26.889	ng/ul	0.00
54) 4-Nitrophenol-d4	14.745	143	35561m	4.260	ng/ul	0.00
60) Fluorene-d10	15.528	176	1100035	29.942	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	219943	26.860	ng/ul	0.00
73) Anthracene-d10	17.386	188	1942496	33.172	ng/ul	0.00
81) Pyrene-d10	19.622	212	2318566	35.917	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.645	264	1960360	34.505	ng/ul	-0.01

Target Compounds Qvalue

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

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