

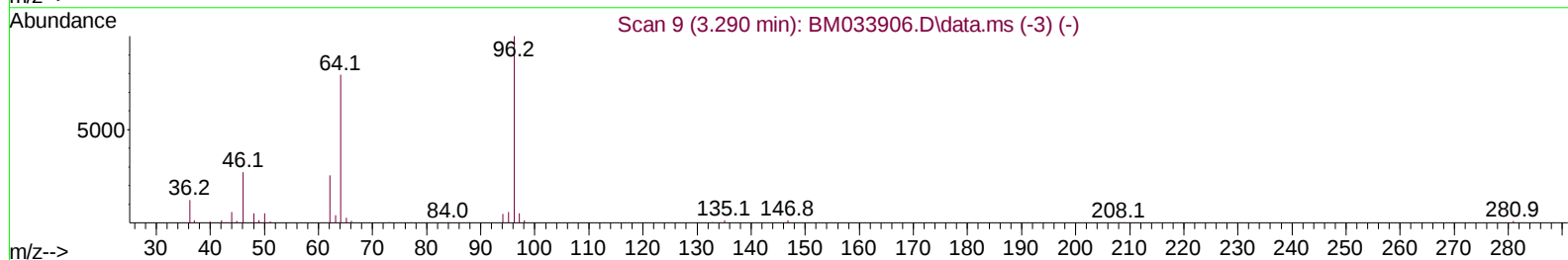
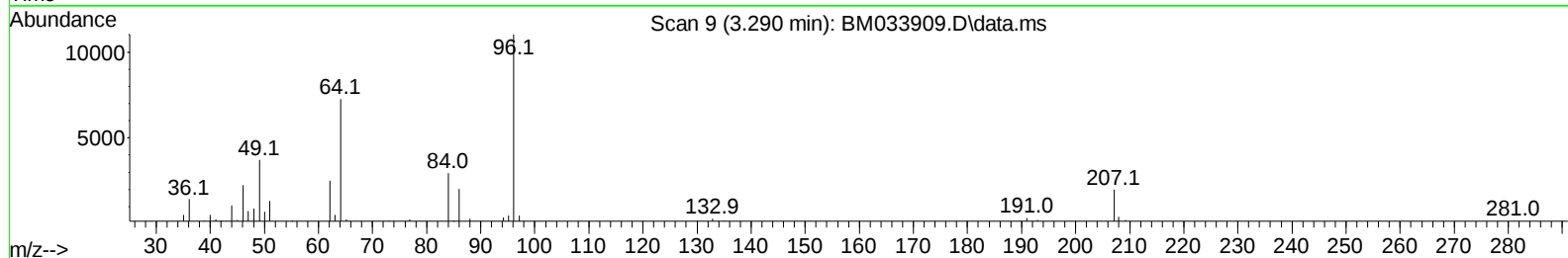
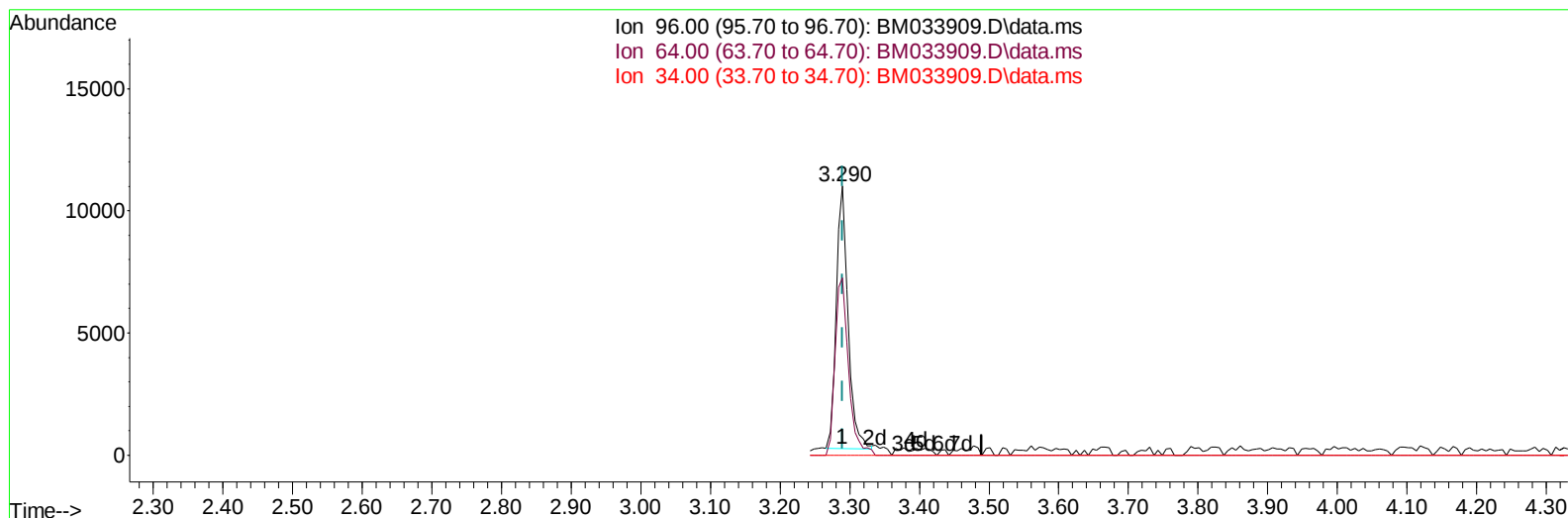
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012822\
 Data File : BM033909.D
 Acq On : 28 Jan 2022 18:20
 Operator : CG/JU
 Sample : N1302-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0F47

Manual IntegrationsAPPROVED

Quant Time: Jan 28 23:28:17 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM010622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jan 28 23:26:29 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/31/2022
 Supervised By :mohammad ahmed 02/01/2022



TIC: BM033909.D\data.ms

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

3.290min (-0.000) 4.11 ng/uL

response 12522

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	65.92
34.00	0.00	0.00
0.00	0.00	0.00

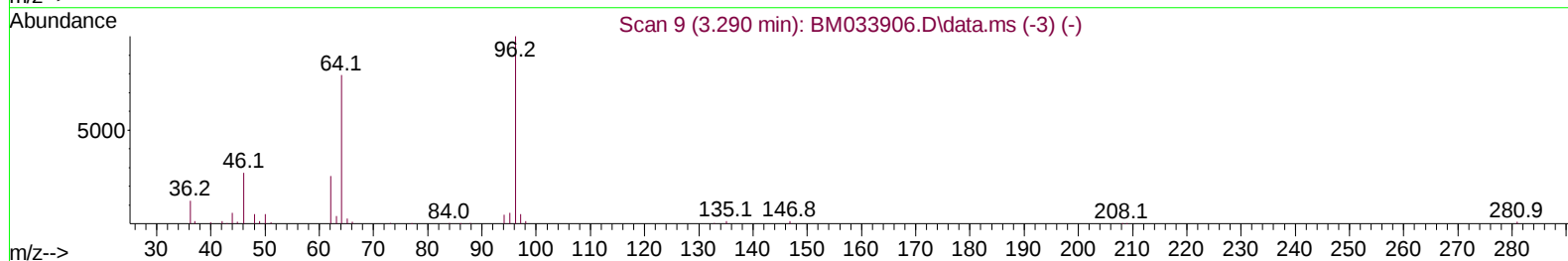
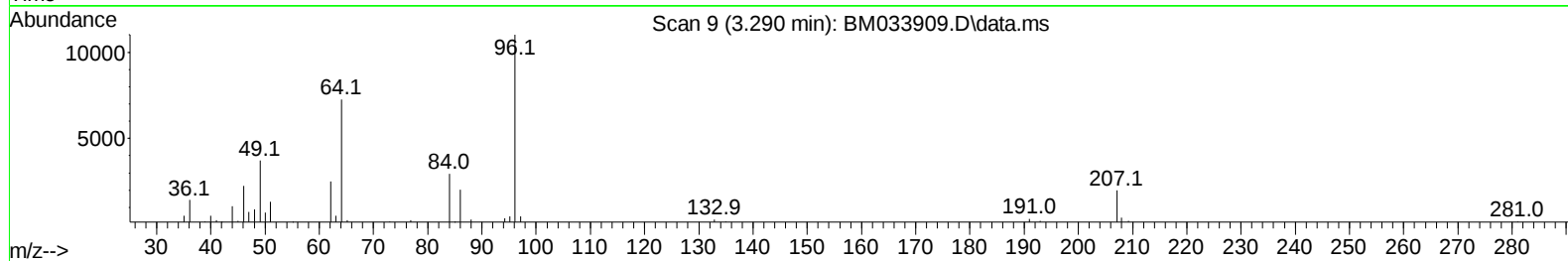
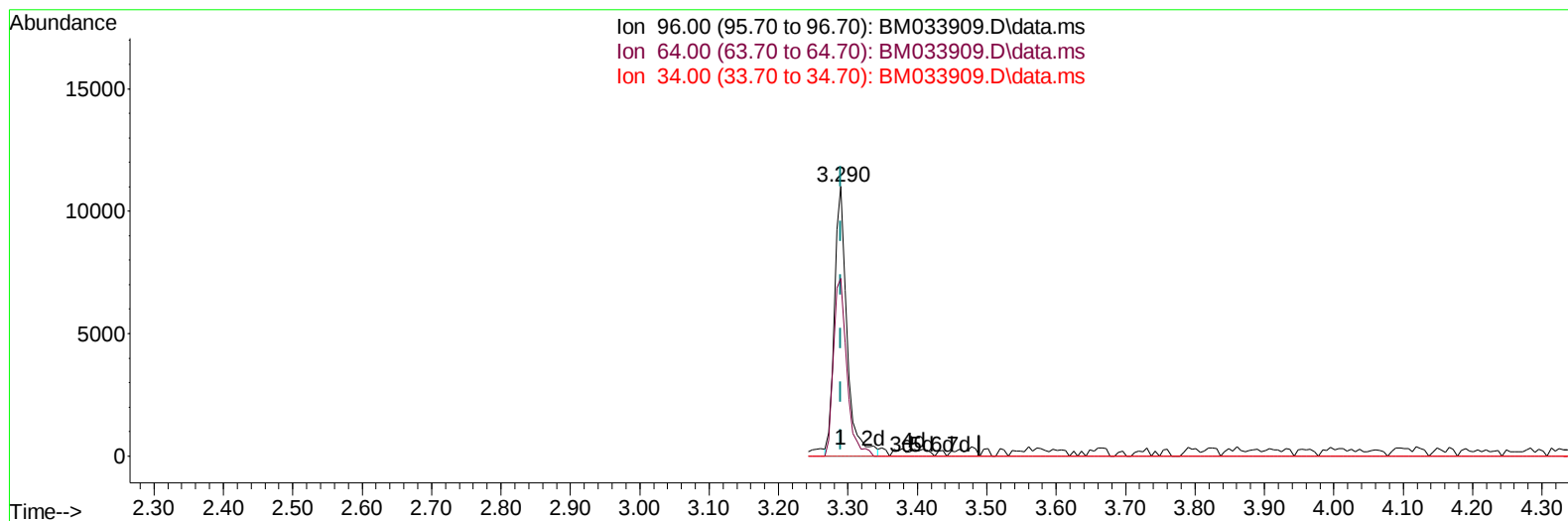
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TIC: BM033909.D\data.ms

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

3.290min (-0.000) 4.51 ng/uL m

response 13763

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	65.92
34.00	0.00	0.00
0.00	0.00	0.00

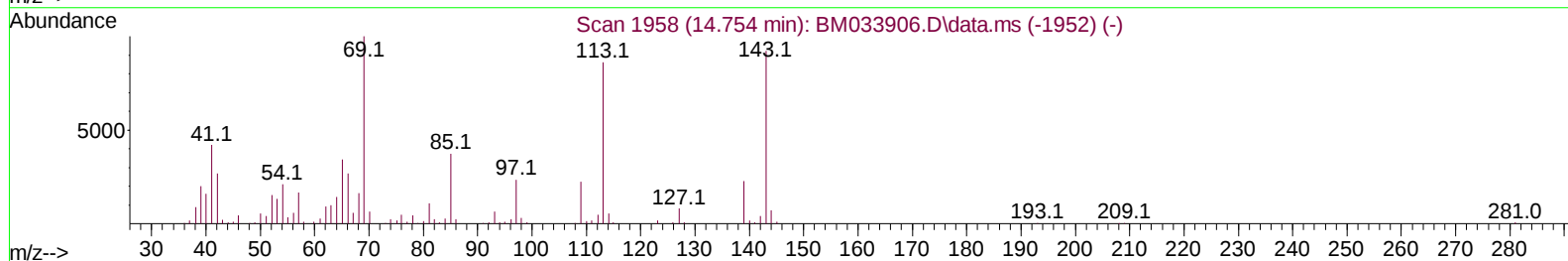
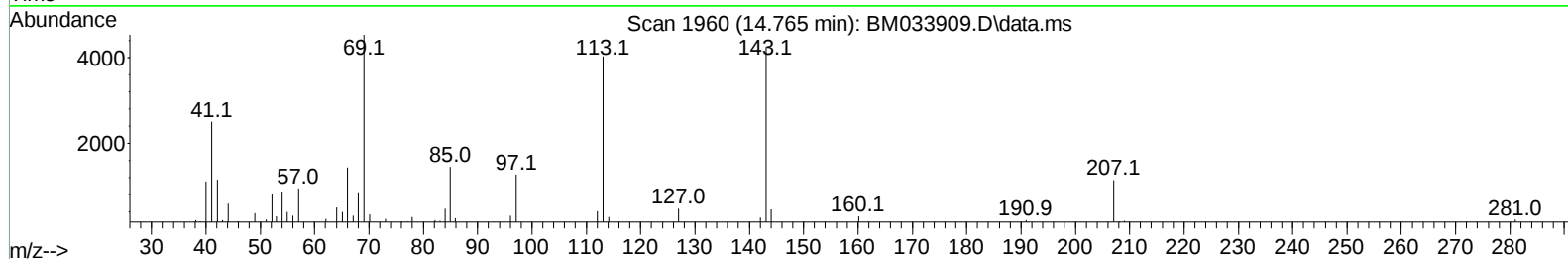
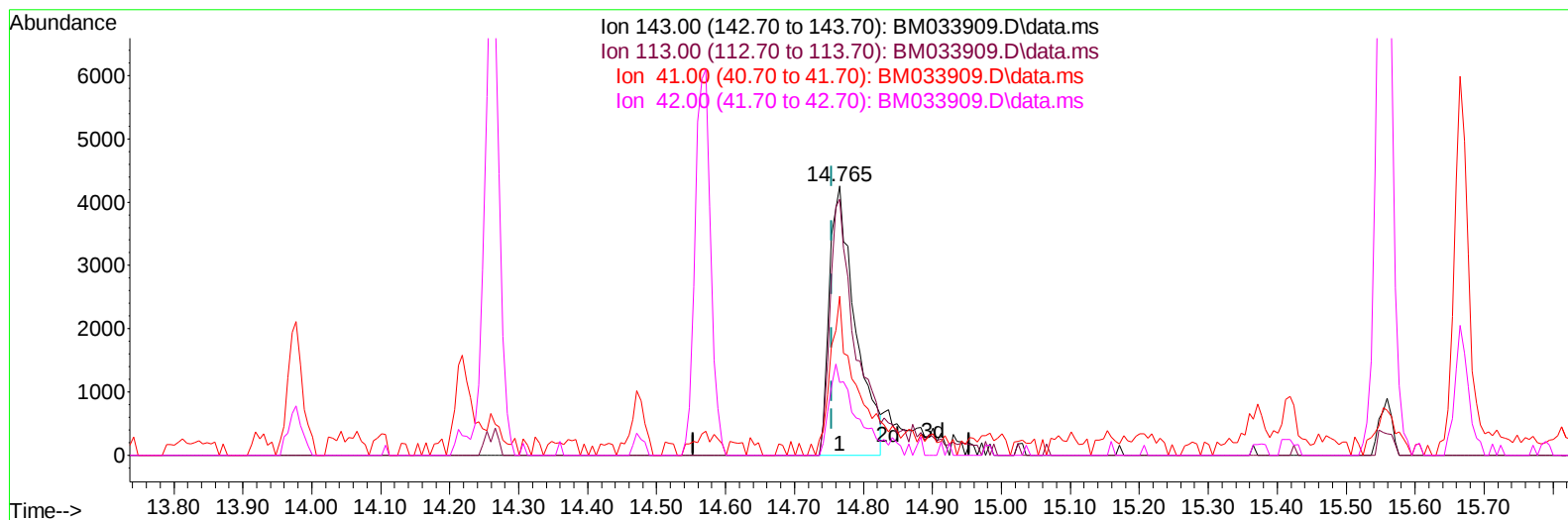
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Operator : CG/JU
Sample : N1302-04
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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TIC: BM033909.D\data.ms

(54) 4-Nitrophenol-d4 (S)**14.765min (+ 0.012) 2.60 ng/u1****response 11056**

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.00	95.04
41.00	57.20	59.06
42.00	39.50	27.09#

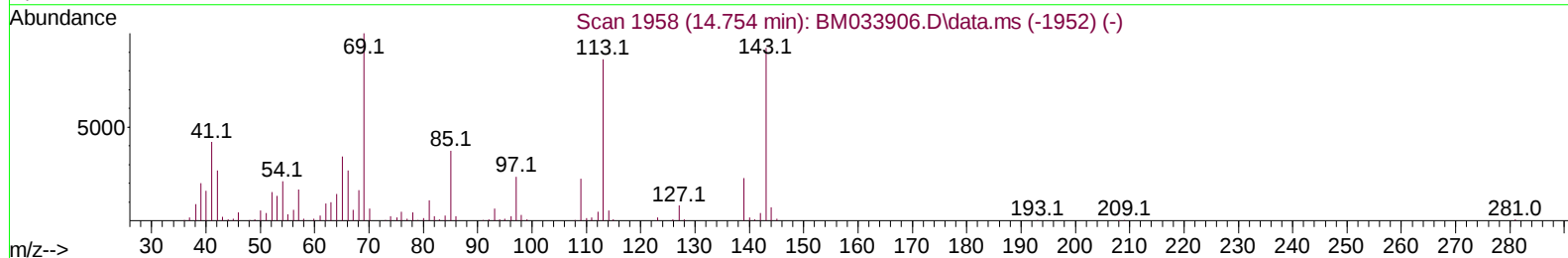
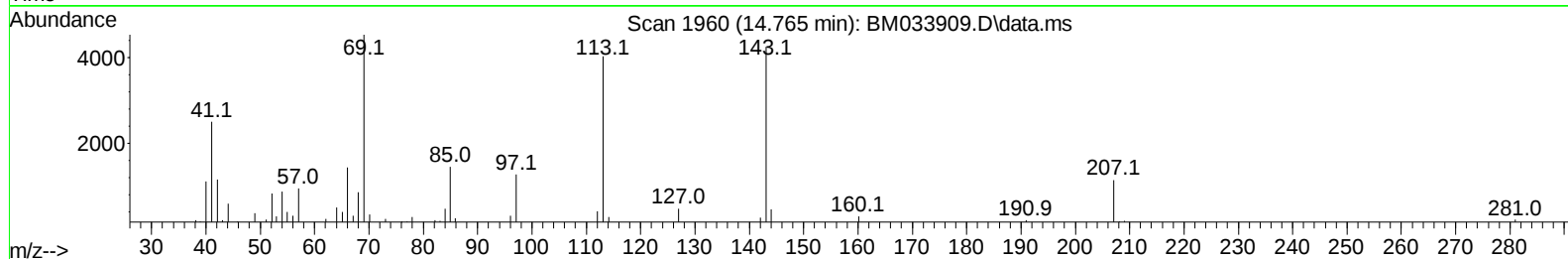
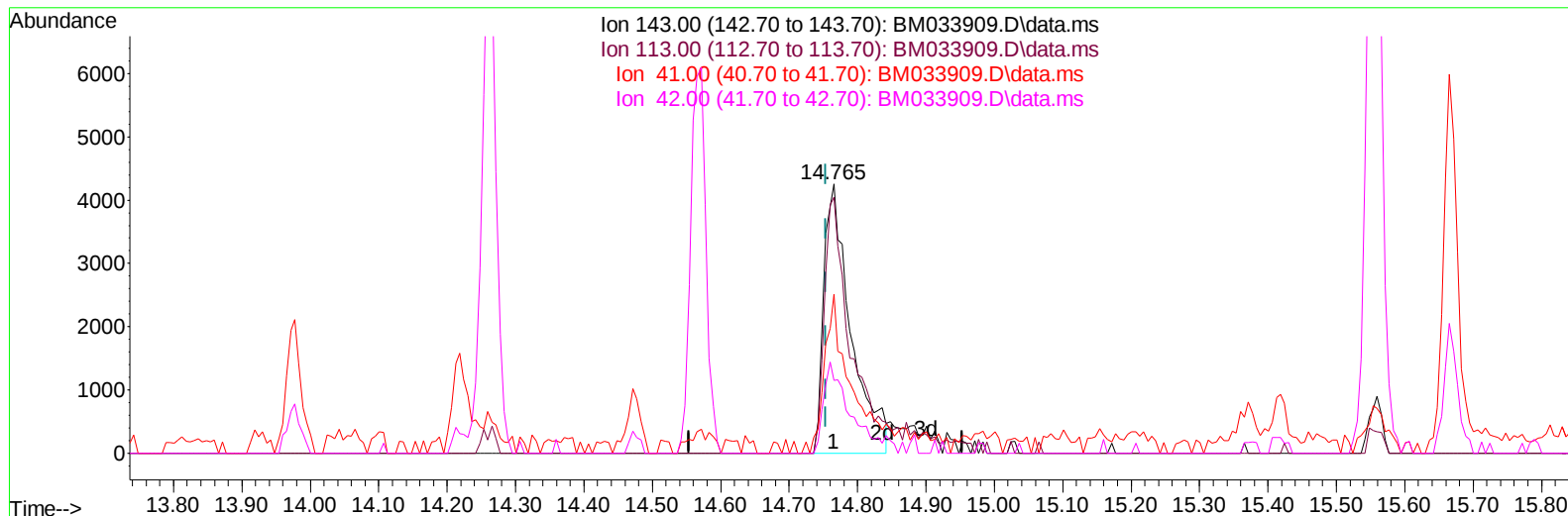
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Operator : CG/JU
Sample : N1302-04
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0F47

Manual IntegrationsAPPROVED

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TIC: BM033909.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.765min (+ 0.012) 2.75 ng/ul m

response 11713

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	105.00	95.04
41.00	57.20	59.06
42.00	39.50	27.09#

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 Misc :
 ALS Vial : 5 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.931	152	114601	20.000	ng/ul	0.00
20) Naphthalene-d8	10.742	136	465872	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.571	164	281577	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.306	188	557460	20.000	ng/ul	0.00
79) Chrysene-d12	21.471	240	523731	20.000	ng/ul	0.00
88) Perylene-d12	23.865	264	524200	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.290	96	13763m	4.512	ng/uL	0.00
4) Pyridine-d5	3.719	84	57107	6.742	ng/ul	0.00
7) Phenol-d5	7.089	99	55760	5.417	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.254	67	191897	30.615	ng/ul	0.00
11) 2-Chlorophenol-d4	7.460	132	185022	23.511	ng/ul	0.00
15) 4-Methylphenol-d8	8.636	113	107666	12.636	ng/ul	0.00
21) Nitrobenzene-d5	9.095	128	118334	32.164	ng/ul	0.00
24) 2-Nitrophenol-d4	9.819	143	119065	30.536	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.360	165	192741	25.784	ng/ul	0.00
31) 4-Chloroaniline-d4	10.877	131	283234	25.490	ng/ul	0.00
46) Dimethylphthalate-d6	13.977	166	716825	32.311	ng/ul	0.00
49) Acenaphthylene-d8	14.265	160	834031	31.867	ng/ul	0.00
54) 4-Nitrophenol-d4	14.765	143	11713m	2.751	ng/ul	0.01
60) Fluorene-d10	15.554	176	612092	33.257	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.671	200	88074	26.560	ng/ul	0.00
73) Anthracene-d10	17.401	188	990916	36.633	ng/ul	0.00
81) Pyrene-d10	19.689	212	1123061	39.708	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.712	264	1045648	38.337	ng/ul	0.00

Target Compounds Qvalue

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

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