

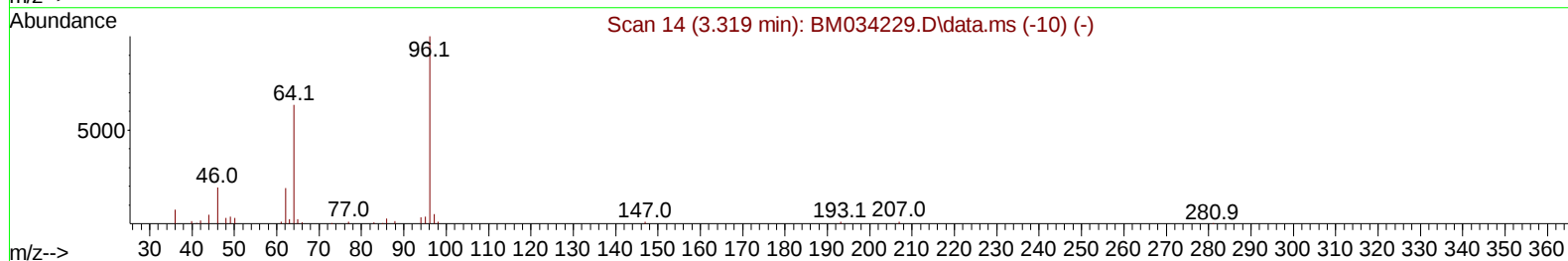
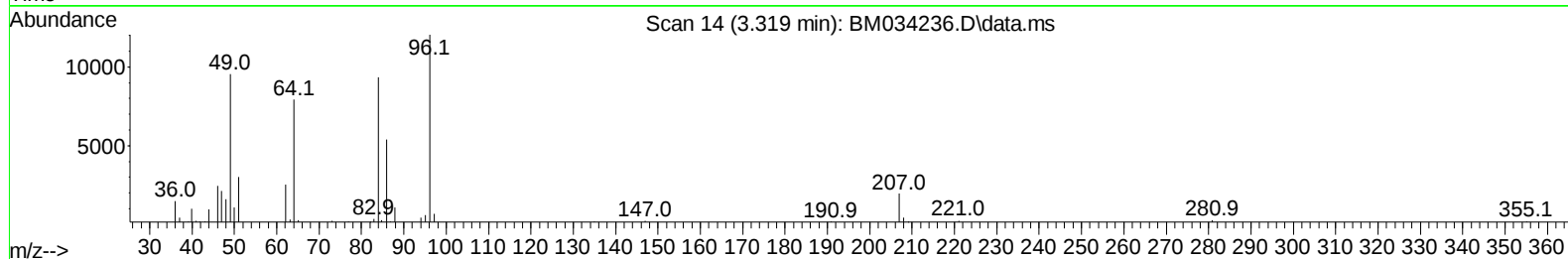
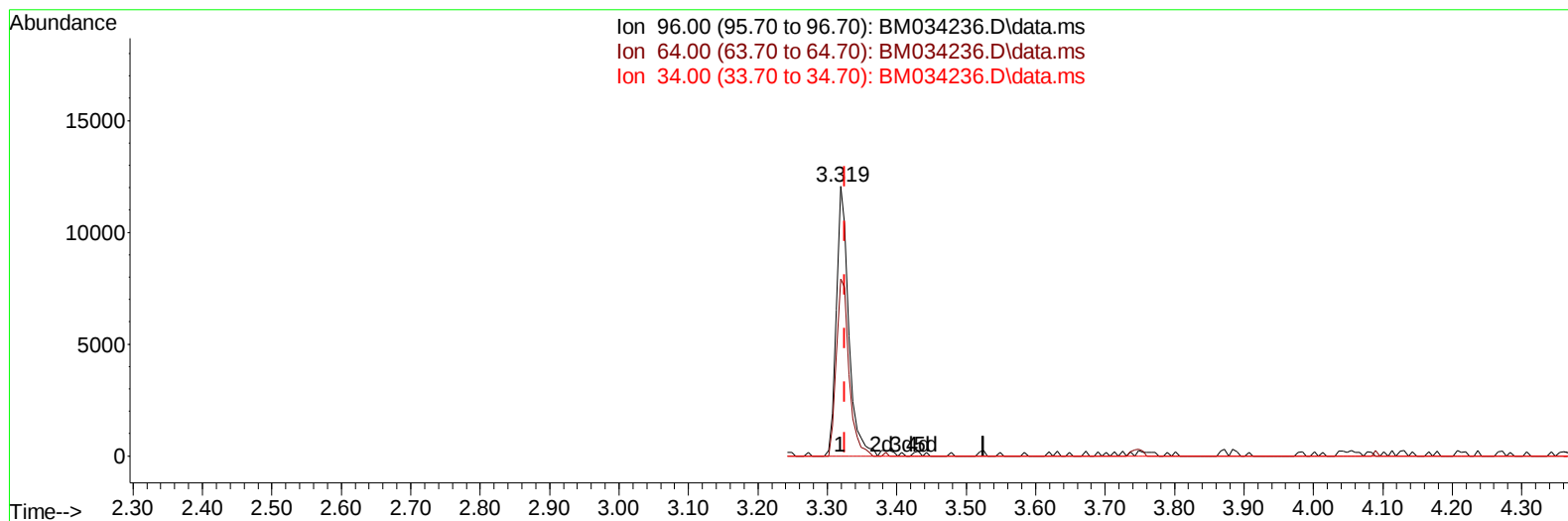
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM031122\
Data File : BM034236.D
Acq On : 12 Mar 2022 00:02
Operator : CG/JU
Sample : N1816-21
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DBSE6

Manual IntegrationsAPPROVED

Quant Time: Mar 12 01:53:05 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM031022.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Mar 11 02:00:05 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/14/2022
Supervised By :mohammad ahmed 03/14/2022



TIC: BM034236.D\data.ms

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

3.319min (-0.006) 4.12 ng/uL

response 14795

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	68.90	65.73
34.00	0.00	0.00
0.00	0.00	0.00

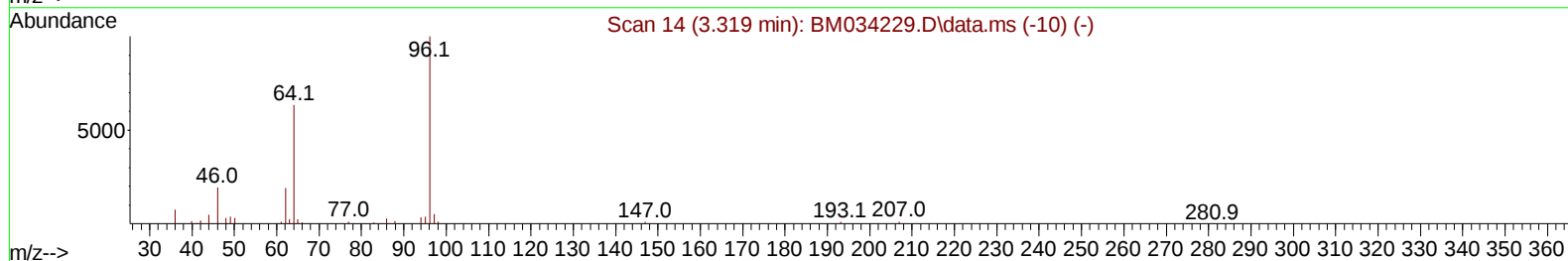
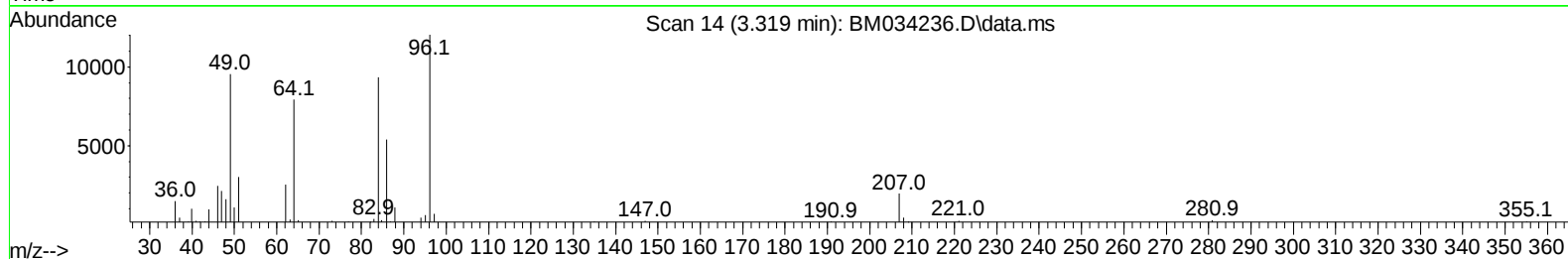
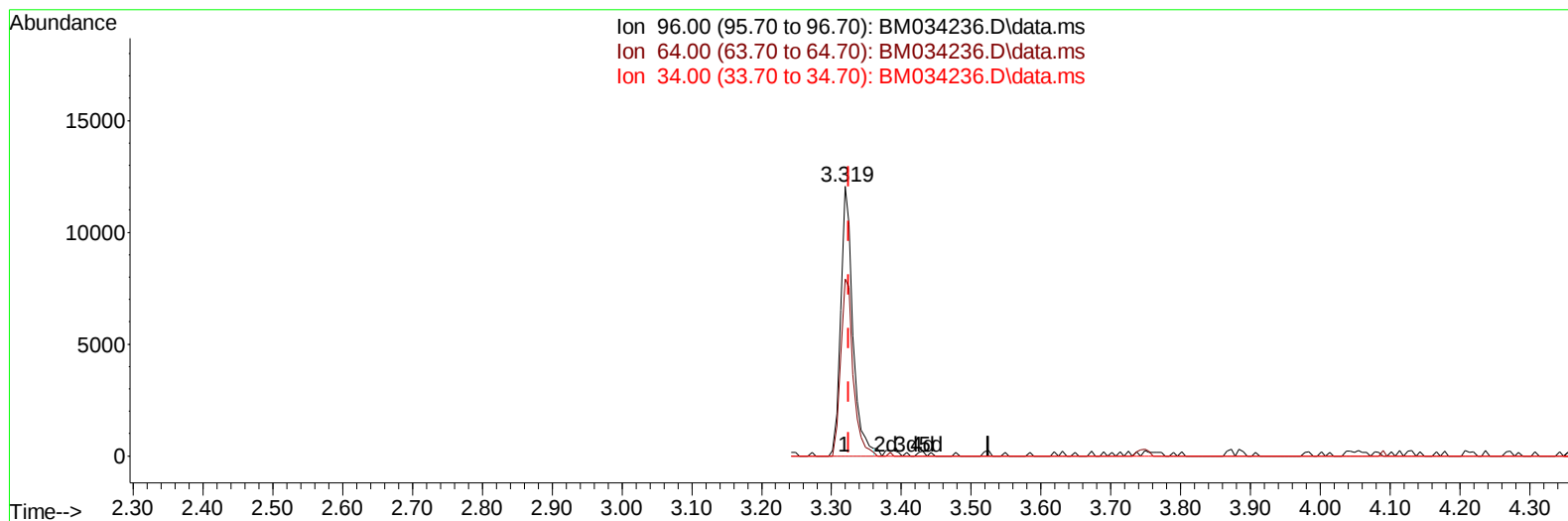
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TIC: BM034236.D\data.ms

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

3.319min (-0.006) 4.08 ng/uL m

response 14677

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	68.90	65.73
34.00	0.00	0.00
0.00	0.00	0.00

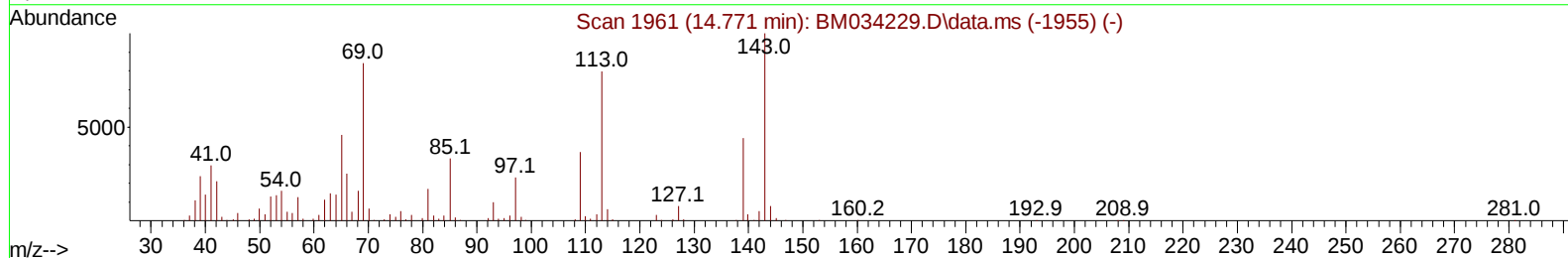
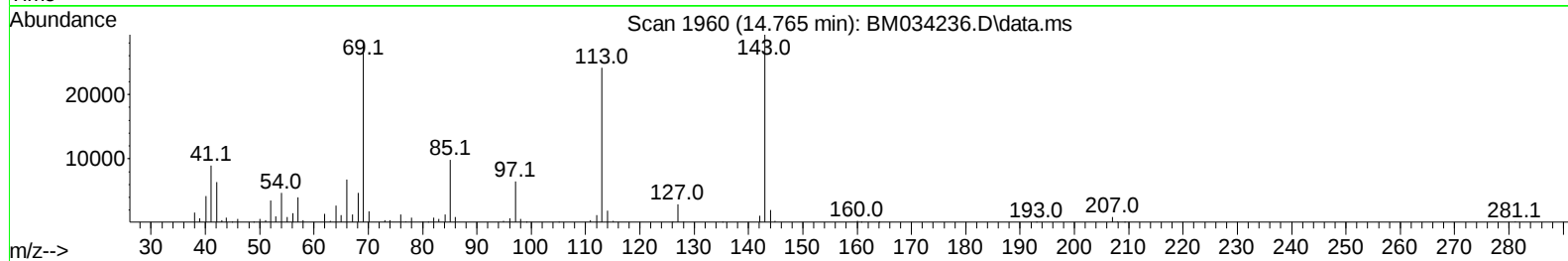
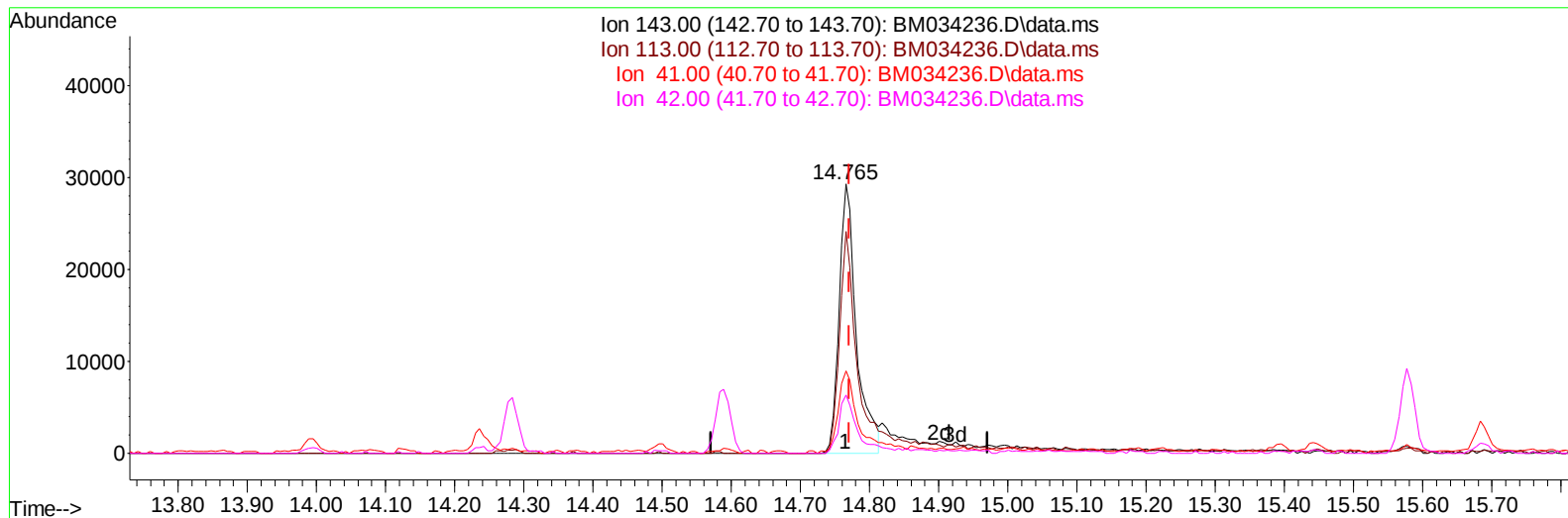
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Instrument :
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TIC: BM034236.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.765min (-0.006) 10.83 ng/u1

response 51018

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	82.35
41.00	31.90	30.53
42.00	19.10	21.64

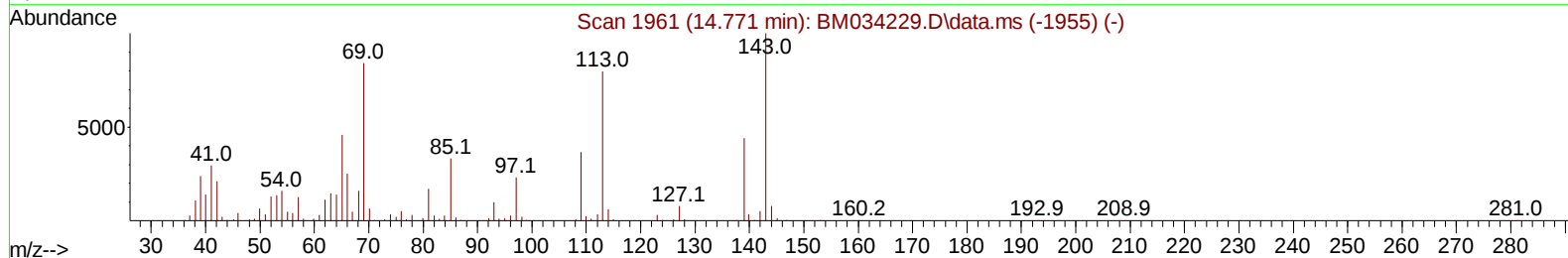
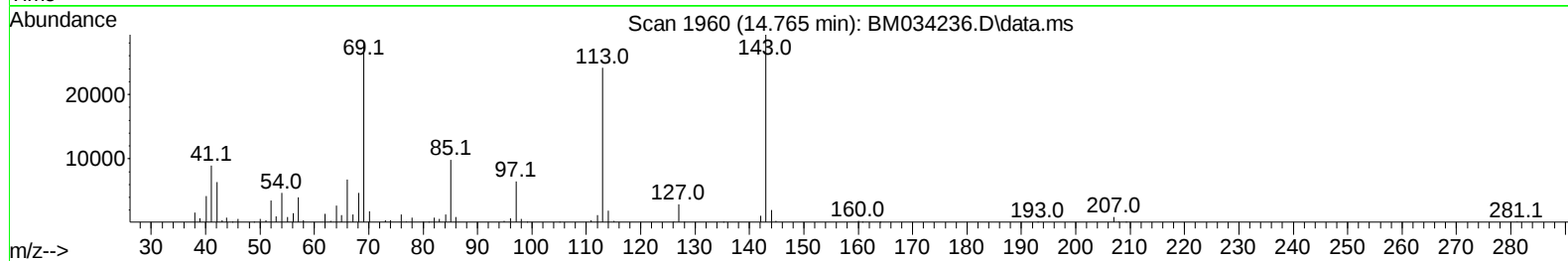
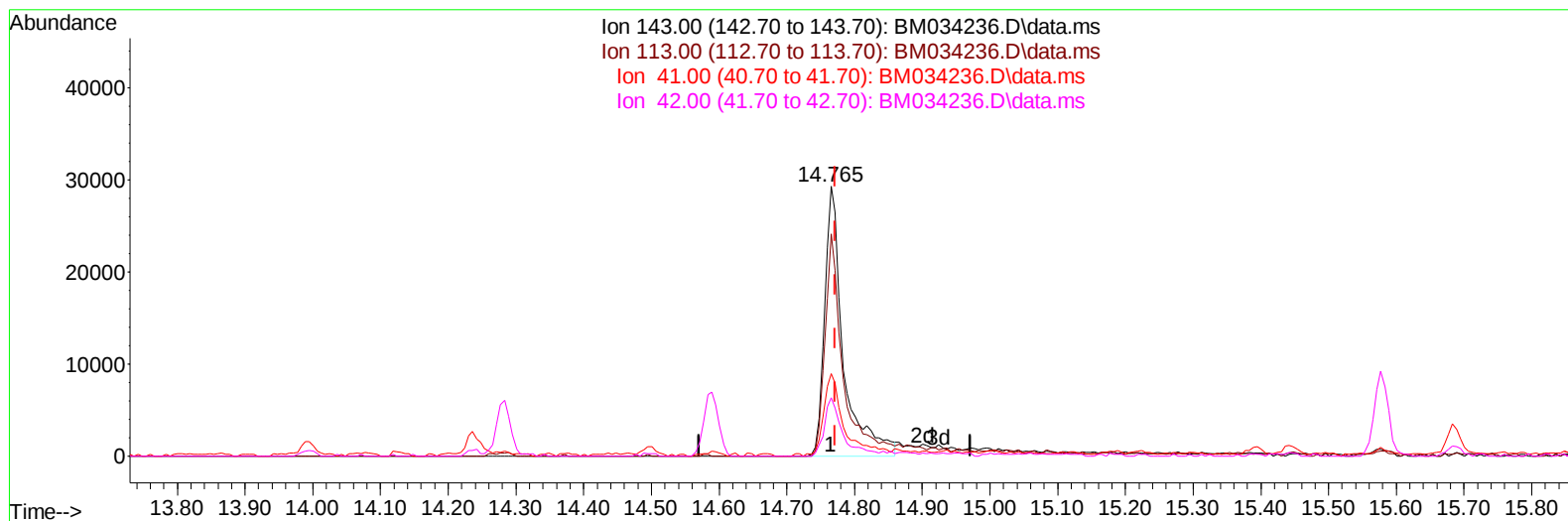
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 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
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TIC: BM034236.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.765min (-0.006) 12.06 ng/ul m

response 56817

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	82.35
41.00	31.90	30.53
42.00	19.10	21.64

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.948	152	156931	20.000	ng/ul	0.00
20) Naphthalene-d8	10.760	136	628444	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.589	164	415604	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.330	188	807740	20.000	ng/ul	0.00
79) Chrysene-d12	21.500	240	695748	20.000	ng/ul	0.00
88) Perylene-d12	23.918	264	711366	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.319	96	14677m	4.082	ng/uL	0.00
4) Pyridine-d5	3.743	84	131341	15.345	ng/ul	0.00
7) Phenol-d5	7.101	99	218441	20.800	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.272	67	145827	23.669	ng/ul	0.00
11) 2-Chlorophenol-d4	7.478	132	207179	22.827	ng/ul	0.00
15) 4-Methylphenol-d8	8.654	113	192049	21.382	ng/ul	0.00
21) Nitrobenzene-d5	9.107	128	103192	22.892	ng/ul	0.00
24) 2-Nitrophenol-d4	9.836	143	103014	20.609	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.378	165	212937	21.250	ng/ul	0.00
31) 4-Chloroaniline-d4	10.889	131	263380	21.135	ng/ul	0.00
46) Dimethylphthalate-d6	13.995	166	706225	23.354	ng/ul	0.00
49) Acenaphthylene-d8	14.283	160	886105	23.071	ng/ul	0.00
54) 4-Nitrophenol-d4	14.765	143	56817m	12.062	ng/ul	0.00
60) Fluorene-d10	15.577	176	627068	23.688	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.689	200	67827	13.272	ng/ul	0.00
73) Anthracene-d10	17.424	188	896204	23.491	ng/ul	0.00
81) Pyrene-d10	19.712	212	1033759	27.292	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.765	264	834715	23.784	ng/ul	0.00
Target Compounds						
					Qvalue	
8) Phenol	7.131	94	22479	2.169	ng/ul	97

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

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