

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102220\
 Data File : BP003711.D
 Acq On : 23 Oct 2020 06:56
 Operator : CG/JU
 Sample : L4221-14
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 OK-03-102020

Quant Time: Oct 23 07:38:31 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP101920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 19 14:54:24 2020
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.528	152	161724	20.00	ng	0.00
21) Naphthalene-d8	11.369	136	607934	20.00	ng	# 0.00
39) Acenaphthene-d10	15.122	164	412077	20.00	ng	0.00
64) Phenanthrene-d10	17.839	188	966706	20.00	ng	0.00
76) Chrysene-d12	21.851	240	1021653	20.00	ng	# 0.00
86) Perylene-d12	24.427	264	928218	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	6.028	112	1205071	124.60	ng	0.00
7) Phenol-d6	7.675	99	1445344	105.78	ng	0.00
23) Nitrobenzene-d5	9.693	82	1217209	84.70	ng	0.00
42) 2,4,6-Tribromophenol	16.598	330	896436	139.44	ng	0.00
45) 2-Fluorobiphenyl	13.751	172	2755785	86.46	ng	0.00
79) Terphenyl-d14	20.310	244	4818022	82.94	ng	0.00
Target Compounds						
10) Phenol	7.699	94	81896	6.29	ng	74
50) Dimethylphthalate	14.540	163	219551	6.48	ng	99

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

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