

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP100419\
 Data File : BP000518.D
 Acq On : 04 Oct 2019 19:37
 Operator : JU
 Sample : K5205-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 RBR-200052-200029

Quant Time: Oct 05 00:37:47 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP100119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 01 18:03:42 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	200019	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	714703	20.00	ng	0.00
39) Acenaphthene-d10	9.81	164	243507	20.00	ng	0.00
64) Phenanthrene-d10	11.31	188	471021	20.00	ng	0.00
76) Chrysene-d12	13.97	240	435467	20.00	ng	0.00
87) Perylene-d12	15.46	264	498227	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	1088200	87.45	ng	0.00
7) Phenol-d6	6.42	99	1339390	89.32	ng	0.00
23) Nitrobenzene-d5	7.33	82	743569	63.54	ng	0.00
42) 2,4,6-Tribromophenol	10.60	330	325980	125.63	ng	0.00
45) 2-Fluorobiphenyl	9.13	172	917213	60.94	ng	0.00
79) Terphenyl-d14	12.92	244	1241861	57.74	ng	0.00
Target Compounds						
10) Phenol	6.42	94	49908	3.251	ng	93
50) Dimethylphthalate	9.52	163	151622	8.961	ng	99

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

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