

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP100719\
 Data File : BP000558.D
 Acq On : 07 Oct 2019 21:21
 Operator : JU
 Sample : K5154-25
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-101I

Quant Time: Oct 08 07:29:37 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	6.76	152	280980	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	1006834	20.00	ng	-0.01
39) Acenaphthene-d10	9.81	164	550923	20.00	ng	0.00
64) Phenanthrene-d10	11.31	188	992533	20.00	ng	0.00
76) Chrysene-d12	13.98	240	879687	20.00	ng	0.00
87) Perylene-d12	15.46	264	872328	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	1311445	75.03	ng	0.00
7) Phenol-d6	6.40	99	1112639	52.82	ng	0.00
23) Nitrobenzene-d5	7.33	82	1291759	78.36	ng	0.00
42) 2,4,6-Tribromophenol	10.61	330	694473	118.29	ng	0.00
45) 2-Fluorobiphenyl	9.13	172	2737282	80.39	ng	0.00
79) Terphenyl-d14	12.92	244	3497216	80.49	ng	0.00
Target Compounds						
2) 1,4-Dioxane	2.47	88	41483	5.943	ng	97
10) Phenol	6.42	94	100062	4.640	ng	94
35) Caprolactam	8.44	113	29188	6.013	ng	94
50) Dimethylphthalate	9.52	163	196444	5.132	ng	99

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

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