

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101119\
 Data File : BP000655.D
 Acq On : 12 Oct 2019 00:47
 Operator : JU
 Sample : K5292-04RE
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 12 03:15:45 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	6.75	152	242494	20.00	ng	-0.02
21) Naphthalene-d8	8.03	136	883413	20.00	ng	-0.02
39) Acenaphthene-d10	9.80	164	487500	20.00	ng	-0.02
64) Phenanthrene-d10	11.29	188	863448	20.00	ng	-0.02
76) Chrysene-d12	13.97	240	716022	20.00	ng	-0.01
87) Perylene-d12	15.45	264	811907	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	204101	13.53	ng	-0.02
7) Phenol-d6	6.39	99	593453	32.65	ng	-0.02
23) Nitrobenzene-d5	7.31	82	840358	58.10	ng	-0.02
42) 2,4,6-Tribromophenol	10.59	330	56149	10.81	ng	-0.02
45) 2-Fluorobiphenyl	9.12	172	1685302	55.93	ng	-0.02
79) Terphenyl-d14	12.90	244	1899567	53.71	ng	-0.01
Target Compounds						
10) Phenol	6.40	94	60206	3.235	ng	96
19) n-Nitroso-di-n-propylamine	7.31	70	103126	12.307	ng	# 73
25) Isophorone	7.42	82	51582	2.009	ng	# 41
41) Hexachlorocyclopentadiene	8.98	237	24	7.301	ng	82
50) Dimethylphthalate	9.51	163	421573	12.445	ng	100
57) 2,4-Dinitrotoluene	9.80	165	63793	6.513	ng	# 32

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

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