

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101119\
 Data File : BP000653.D
 Acq On : 11 Oct 2019 23:58
 Operator : JU
 Sample : K5292-04
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 12 03:15:31 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	240408	20.00	ng	-0.02
21) Naphthalene-d8	8.03	136	889018	20.00	ng	-0.02
39) Acenaphthene-d10	9.80	164	506455	20.00	ng	-0.02
64) Phenanthrene-d10	11.29	188	928825	20.00	ng	-0.02
76) Chrysene-d12	13.96	240	761597	20.00	ng	-0.02
87) Perylene-d12	15.44	264	819485	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	207479	13.87	ng	-0.02
7) Phenol-d6	6.39	99	600872	33.34	ng	-0.02
23) Nitrobenzene-d5	7.31	82	851964	58.53	ng	-0.02
42) 2,4,6-Tribromophenol	10.59	330	58664	10.87	ng	-0.02
45) 2-Fluorobiphenyl	9.12	172	1716659	54.84	ng	-0.02
79) Terphenyl-d14	12.90	244	2085689	55.45	ng	-0.02
Target Compounds						
10) Phenol	6.40	94	61595	3.338	ng	93
19) n-Nitroso-di-n-propylamine	7.31	70	105319	12.678	ng	# 73
25) Isophorone	7.42	82	52455	2.030	ng	# 42
41) Hexachlorocyclopentadiene	9.06	237	11	7.299	ng	93
50) Dimethylphthalate	9.51	163	438423	12.458	ng	99
57) 2,4-Dinitrotoluene	9.80	165	65912	6.478	ng	# 31

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

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