

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP092419\
 Data File : BP000384.D
 Acq On : 24 Sep 2019 15:22
 Operator : JU
 Sample : K4978-04
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 Z2-0163

Quant Time: Sep 24 17:08:35 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP091619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 19 14:44:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	246752	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	869506	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	489790	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	886795	20.00	ng	0.00
76) Chrysene-d12	13.99	240	790905	20.00	ng	-0.01
87) Perylene-d12	15.47	264	875956	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	1246928	87.24	ng	0.00
7) Phenol-d6	6.42	99	1087513	62.07	ng	0.00
23) Nitrobenzene-d5	7.35	82	1299832	96.45	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	730076	150.28	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	2593487	95.30	ng	0.00
79) Terphenyl-d14	12.93	244	3329364	88.31	ng	0.00
Target Compounds						
10) Phenol	6.43	94	62666	3.148	ng	89
27) 2,4-Dimethylphenol	7.74	122	252367	21.350	ng	# 91
50) Dimethylphthalate	9.55	163	451563	13.532	ng	99

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

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