

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101119\
 Data File : BP000648.D
 Acq On : 11 Oct 2019 21:58
 Operator : JU
 Sample : K5292-03RE
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 20190855-AMENDED-COMP-BRE

Quant Time: Oct 12 00:11:45 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	252988	20.00	ng	-0.02
21) Naphthalene-d8	8.03	136	929804	20.00	ng	-0.02
39) Acenaphthene-d10	9.80	164	509170	20.00	ng	-0.02
64) Phenanthrene-d10	11.29	188	923999	20.00	ng	-0.02
76) Chrysene-d12	13.97	240	753651	20.00	ng	-0.01
87) Perylene-d12	15.45	264	826120	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	225360	14.32	ng	-0.02
7) Phenol-d6	6.39	99	644165	33.96	ng	-0.02
23) Nitrobenzene-d5	7.31	82	825481	54.22	ng	-0.02
42) 2,4,6-Tribromophenol	10.59	330	34605	6.38	ng	-0.02
45) 2-Fluorobiphenyl	9.12	172	1574159	50.02	ng	-0.02
79) Terphenyl-d14	12.90	244	1925957	51.74	ng	-0.01
Target Compounds						
10) Phenol	6.40	94	72586	3.738	ng	Qvalue 92
50) Dimethylphthalate	9.51	163	226742	6.409	ng	# 98
75) Fluoranthene	12.52	202	120565	2.313	ng	97
78) Pyrene	12.75	202	161673	3.110	ng	97
81) Benzo(a)anthracene	13.96	228	97487	2.120	ng	# 82
83) Chrysene	13.99	228	90993	2.016	ng	# 90

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

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