

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060916\
 Data File : BF087865.D
 Acq On : 9 Jun 2016 21:19
 Operator : UM/SJ
 Sample : H3378-10RE
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SS-20RE

Quant Time: Jun 10 03:11:21 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 07 18:51:55 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	121624	20.00	ng	-0.02
21) Naphthalene-d8	8.10	136	451814	20.00	ng	-0.02
38) Acenaphthene-d10	9.85	164	195188	20.00	ng	-0.03
63) Phenanthrene-d10	11.32	188	285771	20.00	ng	-0.03
75) Chrysene-d12	13.95	240	249857	20.00	ng	-0.03
86) Perylene-d12	15.36	264	195151	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	379673	53.37	ng	-0.02
7) Phenol-d6	6.44	99	487238	56.51	ng	-0.03
23) Nitrobenzene-d5	7.38	82	289941	37.47	ng	-0.03
41) 2,4,6-Tribromophenol	10.64	330	94468	45.84	ng	-0.03
44) 2-Fluorobiphenyl	9.18	172	459019	33.59	ng	-0.03
78) Terphenyl-d14	12.91	244	342513	31.19	ng	-0.03
Target Compounds						Qvalue
49) Dimethylphthalate	9.58	163	150099	10.62	ng	98
70) Phenanthrene	11.35	178	38930	2.60	ng	97
74) Fluoranthene	12.54	202	75293	4.76	ng	96
77) Pyrene	12.77	202	65759	3.55	ng	97
80) Benzo(a)anthracene	13.94	228	40024	2.81	ng	96
82) Chrysene	13.98	228	30536	2.28	ng	96
87) Benzo(b)fluoranthene	14.96	252	36899m	2.71	ng	
89) Benzo(a)pyrene	15.30	252	26285	2.39	ng	98

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

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