

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121517\
 Data File : BF101468.D
 Acq On : 16 Dec 2017 3:29
 Operator : SJ/JU
 Sample : I6882-05
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 121117-TIDO-E-D-M

Quant Time: Dec 16 05:09:07 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 15 17:36:17 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	150139	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	613162	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	251576	20.00	ng	0.00
63) Phenanthrene-d10	11.35	188	422932	20.00	ng	0.00
75) Chrysene-d12	14.00	240	296394	20.00	ng	0.00
86) Perylene-d12	15.47	264	273453	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	393386	39.03	ng	0.00
7) Phenol-d6	6.45	99	290158	23.13	ng	0.00
23) Nitrobenzene-d5	7.39	82	925787	84.05	ng	0.00
41) 2,4,6-Tribromophenol	10.65	330	200118	82.55	ng	0.00
44) 2-Fluorobiphenyl	9.17	172	1485659	91.61	ng	0.00
78) Terphenyl-d14	12.93	244	1103094	89.72	ng	0.00
Target Compounds						
51) Acenaphthene	9.89	154	85536	5.280	ng	96
54) Dibenzofuran	10.06	168	63233	3.071	ng	99
57) Fluorene	10.40	166	49635	2.850	ng	98
74) Fluoranthene	12.56	202	80522	3.262	ng	99
77) Pyrene	12.80	202	49231	2.213	ng	98

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

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