

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110317\
 Data File : BF100132.D
 Acq On : 3 Nov 2017 14:14
 Operator : SJ/JU
 Sample : I6258-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WOODLAKE-DR-2-14-18

Quant Time: Nov 03 15:51:23 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 03 15:46:10 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	91949	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	384324	20.00	ng	0.00
38) Acenaphthene-d10	9.80	164	182136	20.00	ng	0.00
63) Phenanthrene-d10	11.30	188	335140	20.00	ng	0.00
75) Chrysene-d12	13.95	240	215234	20.00	ng	0.00
86) Perylene-d12	15.41	264	152244	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	517086	88.29	ng	0.00
7) Phenol-d6	6.43	99	616216	88.73	ng	0.00
23) Nitrobenzene-d5	7.34	82	346034	57.97	ng	-0.01
41) 2,4,6-Tribromophenol	10.60	330	151261	91.13	ng	0.00
44) 2-Fluorobiphenyl	9.12	172	702118	59.48	ng	0.00
78) Terphenyl-d14	12.88	244	643827	65.37	ng	0.00
Target Compounds						
10) Phenol	6.44	94	34766	4.560	ng	93
49) Dimethylphthalate	9.52	163	66252	4.593	ng	99
74) Fluoranthene	12.52	202	49990	2.543	ng	99
77) Pyrene	12.75	202	41131	2.513	ng	99

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

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