

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF111117\
 Data File : BF100466.D
 Acq On : 12 Nov 2017 3:57
 Operator : SJ/JU
 Sample : I6408-01
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-A19R(11-13)

Manual Integrations
 APPROVED

SOHIL
 11/13/2017 4:54:00 PM

Quant Time: Nov 13 05:40:09 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 10 15:44:40 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.92	152	58490	20.00	ng	0.02
21) Naphthalene-d8	8.25	136	113173	20.00	ng	0.06
38) Acenaphthene-d10	10.00	164	62727	20.00	ng	0.06
63) Phenanthrene-d10	11.43	188	273661	20.00	ng	0.00
75) Chrysene-d12	14.06	240	298674	20.00	ng	0.00
86) Perylene-d12	15.54	264	203523	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	896286	245.64	ng	0.00
7) Phenol-d6	6.53	99	920093	204.49	ng	0.00
23) Nitrobenzene-d5	7.49	82	335833	196.19	ng	0.02
41) 2,4,6-Tribromophenol	10.76	330	211426	330.77	ng	0.02
44) 2-Fluorobiphenyl	9.33	172	429964	104.37	ng	0.06
78) Terphenyl-d14	13.02	244	1229387	94.58	ng	0.00
Target Compounds						
37) 2-Methylnaphthalene	8.99	142	424153	117.204	ng	# 89
54) Dibenzofuran	10.20	168	63701	11.457	ng	# 23
57) Fluorene	10.53	166	92121m	22.618	ng	
70) Phenanthrene	11.46	178	101880	7.071	ng	# 95

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

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