

Data Path : Z:\HPCHEM1\BNA F\DATA\BF081217\
 Data File : BF097672.D
 Acq On : 13 Aug 2017 13:52
 Operator : SJ/JU
 Sample : I4682-04
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-G10-SSW

Manual Integrations
 APPROVED

sohil
 8/14/2017 7:08:44 PM

Quant Time: Aug 14 08:14:12 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 11 15:55:57 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.50	152	116442	20.00	ng	0.00
21) Naphthalene-d8	9.53	136	514800	20.00	ng	-0.01
38) Acenaphthene-d10	12.35	164	191778	20.00	ng	-0.02
63) Phenanthrene-d10	14.75	188	257270	20.00	ng	-0.01
75) Chrysene-d12	18.45	240	219593	20.00	ng	-0.01
86) Perylene-d12	20.12	264	189349	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	731021	99.79	ng	0.02
7) Phenol-d6	7.07	99	839639	92.39	ng	0.02
23) Nitrobenzene-d5	8.41	82	503094	61.28	ng	-0.01
41) 2,4,6-Tribromophenol	13.65	330	115619	79.31	ng	-0.01
44) 2-Fluorobiphenyl	11.30	172	717650	56.48	ng	-0.02
78) Terphenyl-d14	17.10	244	361232	34.05	ng	-0.02
Target Compounds						
49) Dimethylphthalate	12.00	163	212435	14.373	ng	99
69) Pentachlorophenol	14.52	266	13961	27.671	ng	96
74) Fluoranthene	16.56	202	33458	2.499	ng	94
77) Pyrene	16.86	202	61792	3.403	ng	96
87) Benzo(b)fluoranthene	19.72	252	30998m	2.726	ng	

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

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