

Data Path : Z:\HPCHEM1\BNA F\DATA\BF081117\
 Data File : BF097616.D
 Acq On : 12 Aug 2017 2:47
 Operator : SJ/JU
 Sample : I4631-07
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VE-3-1

Quant Time: Aug 14 11:01:01 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	7.51	152	92213	20.00	ng	0.00
21) Naphthalene-d8	9.54	136	327757	20.00	ng	0.00
38) Acenaphthene-d10	12.37	164	122387	20.00	ng	0.00
63) Phenanthrene-d10	14.79	188	148515	20.00	ng	0.03
75) Chrysene-d12	18.50	240	179891	20.00	ng	0.04
86) Perylene-d12	20.16	264	169156	20.00	ng	0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	398452	68.68	ng	0.04
7) Phenol-d6	7.10	99	277654	38.58	ng	0.04
23) Nitrobenzene-d5	8.43	82	476020	91.07	ng	0.00
41) 2,4,6-Tribromophenol	13.70	330	109443	117.63	ng	0.04
44) 2-Fluorobiphenyl	11.33	172	675282	83.27	ng	0.00
78) Terphenyl-d14	17.16	244	470033	54.08	ng	0.04
Target Compounds						
10) Phenol	7.13	94	46960	5.960	ng	82
17) 2-Methylphenol	8.06	107	63902	12.398	ng	# 95
20) 3+4-Methylphenols	8.31	107	372388	58.993	ng	# 57
27) 2,4-Dimethylphenol	9.13	122	78230	16.282	ng	98
31) Naphthalene	9.58	128	41420	2.523	ng	# 84
37) 2-Methylnaphthalene	10.71	142	339982	32.914	ng	97
51) Acenaphthene	12.43	154	44121	5.744	ng	94
54) Dibenzofuran	12.72	168	39799	3.701	ng	# 77
57) Fluorene	13.27	166	59638	6.744	ng	# 97
70) Phenanthrene	14.83	178	94747	11.077	ng	# 88

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

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