

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF053017\
 Data File : BF095555.D
 Acq On : 31 May 2017 5:03
 Operator : SJ/MA
 Sample : I3388-06
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-B29-B35-002

Quant Time: May 31 08:45:17 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 26 16:50:16 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	79539	20.00	ng	0.00
21) Naphthalene-d8	9.75	136	302755	20.00	ng	0.00
38) Acenaphthene-d10	12.60	164	119631	20.00	ng	0.00
63) Phenanthrene-d10	15.01	188	180106	20.00	ng	0.00
75) Chrysene-d12	18.68	240	176572	20.00	ng	0.00
86) Perylene-d12	20.37	264	100951	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.72	112	622958	130.58	ng	0.01
7) Phenol-d6	7.30	99	675486	118.01	ng	0.00
23) Nitrobenzene-d5	8.64	82	433277	90.24	ng	-0.01
41) 2,4,6-Tribromophenol	13.91	330	120287	122.77	ng	0.00
44) 2-Fluorobiphenyl	11.53	172	846534	101.85	ng	0.00
78) Terphenyl-d14	17.32	244	654522	81.65	ng	0.00
Target Compounds						Qvalue
17) 2-Methylphenol	8.35	107	11769	2.65	ng	# 95
27) 2,4-Dimethylphenol	9.38	122	40609	9.25	ng	97
31) Naphthalene	9.79	128	1256785	82.59	ng	100
37) 2-Methylnaphthalene	10.92	142	112960	11.31	ng	96
51) Acenaphthene	12.65	154	25419	3.34	ng	96
57) Fluorene	13.50	166	18531	2.02	ng	# 85
70) Phenanthrene	15.05	178	23771	2.30	ng	97
72) Carbazole	15.44	167	52752	5.48	ng	97

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

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