

Data Path : Z:\HPCHEM1\BNA F\DATA\BF081217\
 Data File : BF097646.D
 Acq On : 12 Aug 2017 21:27
 Operator : SJ/JU
 Sample : I4631-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VE-3-1

Quant Time: Aug 14 12:00:45 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	122099	20.00	ng	0.00
21) Naphthalene-d8	9.53	136	464806	20.00	ng	0.00
38) Acenaphthene-d10	12.37	164	150816	20.00	ng	0.00
63) Phenanthrene-d10	14.78	188	178793	20.00	ng	0.02
75) Chrysene-d12	18.49	240	177074	20.00	ng	0.02
86) Perylene-d12	20.14	264	166235	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	484341	63.05	ng	0.02
7) Phenol-d6	7.09	99	393051	41.25	ng	0.04
23) Nitrobenzene-d5	8.42	82	661214	89.20	ng	0.00
41) 2,4,6-Tribromophenol	13.69	330	131225	114.46	ng	0.03
44) 2-Fluorobiphenyl	11.32	172	950956	95.16	ng	0.00
78) Terphenyl-d14	17.14	244	473206	55.31	ng	0.02
Target Compounds						
10) Phenol	7.12	94	68112	6.528	ng	94
14) 1,2-Dichlorobenzene	7.76	146	20457	2.233	ng	91
17) 2-Methylphenol	8.05	107	81029	11.873	ng	95
20) 3+4-Methylphenols	8.30	107	528114	63.185	ng	# 58
27) 2,4-Dimethylphenol	9.13	122	125002	18.346	ng	98
31) Naphthalene	9.57	128	60003	2.577	ng	# 81
37) 2-Methylnaphthalene	10.70	142	435627	29.738	ng	97
51) Acenaphthene	12.42	154	54783	5.788	ng	91
54) Dibenzofuran	12.71	168	55755	4.208	ng	# 70
57) Fluorene	13.26	166	83730	7.684	ng	# 96
70) Phenanthrene	14.82	178	112895	10.963	ng	# 90

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

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