

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF121417\
 Data File : BF101419.D
 Acq On : 15 Dec 2017 2:57
 Operator : SJ/JU
 Sample : I6824-04
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TW-3

Quant Time: Dec 15 13:59:01 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 14 15:42:32 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	174709	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	696165	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	320531	20.00	ng	0.00
63) Phenanthrene-d10	11.35	188	588825	20.00	ng	0.00
75) Chrysene-d12	13.99	240	418014	20.00	ng	-0.01
86) Perylene-d12	15.46	264	338171	20.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.43	112	678073	57.81	ng	0.00
7) Phenol-d6	6.45	99	550395	37.71	ng	-0.01
23) Nitrobenzene-d5	7.39	82	1087098	86.93	ng	0.00
41) 2,4,6-Tribromophenol	10.65	330	338428	109.57	ng	0.00
44) 2-Fluorobiphenyl	9.17	172	1723291	83.40	ng	0.00
78) Terphenyl-d14	12.93	244	1640689	94.62	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
10) Phenol	6.46	94	61949	3.724	ng	94
17) 2-Methylphenol	7.07	107	80366	7.081	ng	99
20) 3+4-Methylphenols	7.23	107	226190	18.808	ng	93
27) 2,4-Dimethylphenol	7.75	122	234068	23.170	ng	98
31) Naphthalene	8.12	128	2268400	64.724	ng	99
37) 2-Methylnaphthalene	8.81	142	243691	10.672	ng	99
45) 1,1'-Biphenyl	9.27	154	59508	2.093	ng	97
51) Acenaphthene	9.88	154	194998	9.448	ng	100
54) Dibenzofuran	10.06	168	202083	7.704	ng	98
57) Fluorene	10.40	166	166810	7.518	ng	97
70) Phenanthrene	11.37	178	280616	8.570	ng	99
71) Anthracene	11.42	178	73661	2.202	ng	97
72) Carbazole	11.58	167	401191	12.811	ng	98

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

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