

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF121417\
 Data File : BF101429.D
 Acq On : 15 Dec 2017 7:51
 Operator : SJ/JU
 Sample : I6878-02
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TW-3-DEP

Manual Integrations
 APPROVED

Sohil
 12/15/2017 6:45:15 PM

Quant Time: Dec 15 10:29:44 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	6.82	152	170819	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	659058	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	287547	20.00	ng	0.00
63) Phenanthrene-d10	11.35	188	485865	20.00	ng	0.00
75) Chrysene-d12	13.99	240	334628	20.00	ng	-0.01
86) Perylene-d12	15.47	264	324585	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	395885	34.52	ng	0.00
7) Phenol-d6	6.45	99	299772	21.00	ng	0.00
23) Nitrobenzene-d5	7.39	82	913237	77.13	ng	0.00
41) 2,4,6-Tribromophenol	10.65	330	198808	71.75	ng	0.00
44) 2-Fluorobiphenyl	9.17	172	1507562	81.33	ng	0.00
78) Terphenyl-d14	12.93	244	1111903	80.11	ng	0.00
Target Compounds						
10) Phenol	6.47	94	34031	2.092	ng	79
17) 2-Methylphenol	7.07	107	67422	6.076	ng	96
20) 3+4-Methylphenols	7.23	107	163848	13.934	ng	89
27) 2,4-Dimethylphenol	7.75	122	191584	20.032	ng	98
31) Naphthalene	8.12	128	988748	29.800	ng	99
37) 2-Methylnaphthalene	8.81	142	127726	5.909	ng	99
51) Acenaphthene	9.89	154	171142	9.243	ng	99
54) Dibenzofuran	10.06	168	184485	7.840	ng	98
57) Fluorene	10.40	166	143752	7.222	ng	97
70) Phenanthrene	11.37	178	282477	10.454	ng	99
71) Anthracene	11.42	178	70193	2.543	ng	97
72) Carbazole	11.58	167	328175	12.700	ng	99

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

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