

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122617\
 Data File : BF101709.D
 Acq On : 26 Dec 2017 12:24
 Operator : SJ/JU
 Sample : I6835-03DL 5X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-4539-02DL

Manual Integrations
 APPROVED

Sohil
 12/27/2017 4:14:19 PM

Quant Time: Dec 26 13:14:15 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 26 10:48:01 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	152453	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	601079	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	290918	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	578441	20.00	ng	0.00
75) Chrysene-d12	13.96	240	529436	20.00	ng	0.00
86) Perylene-d12	15.43	264	464504	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	292578	28.59	ng	0.01
7) Phenol-d6	6.43	99	350295	27.50	ng	0.00
23) Nitrobenzene-d5	7.36	82	249567	23.11	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	100756	35.94	ng	0.00
44) 2-Fluorobiphenyl	9.14	172	483530	25.78	ng	0.00
78) Terphenyl-d14	12.89	244	576757	26.26	ng	0.00
Target Compounds						
10) Phenol	6.45	94	331979	22.867	ng	99
17) 2-Methylphenol	7.05	107	162007	16.359	ng	99
20) 3+4-Methylphenols	7.21	107	403200	38.421	ng	84
27) 2,4-Dimethylphenol	7.73	122	160881	18.445	ng	99
31) Naphthalene	8.09	128	2508960	82.912	ng	98
37) 2-Methylnaphthalene	8.79	142	117657	5.968	ng	99
48) Acenaphthylene	9.69	152	185749	5.957	ng	99
54) Dibenzofuran	10.03	168	75686	3.179	ng	98
57) Fluorene	10.37	166	71621	3.556	ng	97
70) Phenanthrene	11.34	178	95078	2.956	ng	98
72) Carbazole	11.56	167	234964	7.638	ng	98

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

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