

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF012318\
 Data File : BF102380.D
 Acq On : 24 Jan 2018 9:44
 Operator : SJ/JU
 Sample : J1239-06DL 10X
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-14R30-02DL

Manual Integrations
 APPROVED

Sohil
 1/24/2018 2:47:37 PM

Quant Time: Jan 24 13:57:46 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 23 00:43:41 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	121601	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	508920	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	214762	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	312851	20.00	ng	0.00
75) Chrysene-d12	13.87	240	238582	20.00	ng	0.00
86) Perylene-d12	15.29	264	210110	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	110221	13.74	ng	0.00
7) Phenol-d6	6.35	99	117395m	12.14	ng	0.00
23) Nitrobenzene-d5	7.27	82	87851	9.92	ng	-0.01
41) 2,4,6-Tribromophenol	10.54	330	24931	11.72	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	185866	12.73	ng	-0.01
78) Terphenyl-d14	12.82	244	131307	12.41	ng	0.00
Target Compounds						
10) Phenol	6.37	94	487356	46.171	ng	99
17) 2-Methylphenol	6.97	107	116801	15.997	ng	99
20) 3+4-Methylphenols	7.13	107	349855	40.742	ng	84
27) 2,4-Dimethylphenol	7.66	122	67514	9.748	ng	92
31) Naphthalene	8.02	128	1251682	49.260	ng	100
37) 2-Methylnaphthalene	8.71	142	36517	2.158	ng	96
48) Acenaphthylene	9.61	152	85651	3.683	ng	98
72) Carbazole	11.47	167	97831	5.389	ng	98

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

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