

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF012318\
 Data File : BF102379.D
 Acq On : 24 Jan 2018 9:17
 Operator : SJ/JU
 Sample : J1239-13DL 25X
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-11R-01DL

Quant Time: Jan 24 10:46:22 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	125480	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	515034	20.00	ng	0.00
38) Acenaphthene-d10	9.74	164	220090	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	335172	20.00	ng	0.00
75) Chrysene-d12	13.87	240	249064	20.00	ng	0.00
86) Perylene-d12	15.29	264	221905	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	38239	4.62	ng	0.00
7) Phenol-d6	6.35	99	46060	4.62	ng	0.00
23) Nitrobenzene-d5	7.27	82	32554	3.63	ng	-0.01
41) 2,4,6-Tribromophenol	10.54	330	8772	4.02	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	74321	4.97	ng	-0.01
78) Terphenyl-d14	12.82	244	53812	4.87	ng	0.00
Target Compounds						
3) Pyridine	3.10	79	26579	2.587	ng	Qvalue 97
10) Phenol	6.37	94	1012759	92.981	ng	98
17) 2-Methylphenol	6.97	107	169338	22.476	ng	95
20) 3+4-Methylphenols	7.13	107	565790	63.852	ng	87
27) 2,4-Dimethylphenol	7.66	122	81954	11.692	ng	97
31) Naphthalene	8.02	128	894051	34.768	ng	100
48) Acenaphthylene	9.61	152	54126	2.271	ng	98
72) Carbazole	11.47	167	52280	2.688	ng	97
76) Benzydine	12.57	184	28372	3.391	ng	# 94

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

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