

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF012318\
 Data File : BF102352.D
 Acq On : 23 Jan 2018 19:57
 Operator : SJ/JU
 Sample : J1015-14
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-01-012218-B

Manual Integrations
 APPROVED

Sohil
 1/24/2018 2:46:58 PM

Quant Time: Jan 24 02:26:49 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	151440	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	642320	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	287995	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	499742	20.00	ng	0.00
75) Chrysene-d12	13.87	240	331041	20.00	ng	0.00
86) Perylene-d12	15.30	264	299455	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	958780	96.00	ng	0.00
7) Phenol-d6	6.36	99	1165189	96.77	ng	0.00
23) Nitrobenzene-d5	7.28	82	716969	64.15	ng	0.00
41) 2,4,6-Tribromophenol	10.54	330	242225	84.88	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	1175495	60.01	ng	0.00
78) Terphenyl-d14	12.82	244	1018209	69.34	ng	0.00
Target Compounds						
10) Phenol	6.37	94	52209	3.972	ng	97
49) Dimethylphthalate	9.47	163	120640	5.068	ng	99
70) Phenanthrene	11.26	178	216217	7.425	ng	97
74) Fluoranthene	12.44	202	228144	7.683	ng	98
77) Pyrene	12.67	202	311815	12.183	ng	99
80) Benzo(a)anthracene	13.86	228	90680	4.449	ng	93
82) Chrysene	13.90	228	112892	5.455	ng	96
85) Indeno(1,2,3-cd)pyrene	16.67	276	41691	2.439	ng	98
87) Benzo(b)fluoranthene	14.89	252	83885m	4.322	ng	
89) Benzo(a)pyrene	15.23	252	67886	3.878	ng	# 92
91) Benzo(g,h,i)perylene	17.09	276	49200	3.514	ng	# 90

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

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