

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF012318\
 Data File : BF102353.D
 Acq On : 23 Jan 2018 20:24
 Operator : SJ/JU
 Sample : J1015-16
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 1/24/2018 2:47:01 PM

Quant Time: Jan 24 02:32:04 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	6.71	152	134190	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	567757	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	256725	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	441073	20.00	ng	0.00
75) Chrysene-d12	13.87	240	280206	20.00	ng	0.00
86) Perylene-d12	15.30	264	257535	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	868284	98.11	ng	0.00
7) Phenol-d6	6.36	99	1050136	98.42	ng	0.00
23) Nitrobenzene-d5	7.28	82	652185	66.01	ng	0.00
41) 2,4,6-Tribromophenol	10.54	330	216959	85.29	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	987410	56.55	ng	0.00
78) Terphenyl-d14	12.82	244	786603	63.29	ng	0.00
Target Compounds						
10) Phenol	6.37	94	53307	4.576	ng	Qvalue 97
49) Dimethylphthalate	9.47	163	152514	7.187	ng	100
70) Phenanthrene	11.26	178	110468	4.298	ng	98
74) Fluoranthene	12.45	202	192040	7.327	ng	97
77) Pyrene	12.67	202	183336	8.462	ng	98
80) Benzo(a)anthracene	13.86	228	68700	3.982	ng	98
82) Chrysene	13.89	228	75291	4.298	ng	97
85) Indeno(1,2,3-cd)pyrene	16.67	276	35824	2.476	ng	# 94
87) Benzo(b)fluoranthene	14.89	252	83058m	4.976	ng	
89) Benzo(a)pyrene	15.23	252	56912	3.780	ng	# 92
91) Benzo(g,h,i)perylene	17.10	276	36044	2.993	ng	99

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

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