

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF021218\
 Data File : BF102913.D
 Acq On : 12 Feb 2018 11:56
 Operator : SJ/JU
 Sample : J1512-10
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RO-4632

Manual Integrations
 APPROVED

Sohil
 2/13/2018 11:25:14 AM

Quant Time: Feb 12 12:56:11 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 05 12:44:16 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	187075	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	781319	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	346779	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	556365	20.00	ng	0.00
75) Chrysene-d12	14.05	240	353383	20.00	ng	0.00
86) Perylene-d12	15.54	264	359132	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1200858	97.49	ng	0.01
7) Phenol-d6	6.52	99	1425617	96.12	ng	0.01
23) Nitrobenzene-d5	7.45	82	929672	82.54	ng	0.00
41) 2,4,6-Tribromophenol	10.72	330	269650	96.61	ng	0.00
44) 2-Fluorobiphenyl	9.24	172	1396437	66.82	ng	0.00
78) Terphenyl-d14	13.00	244	1024618	68.65	ng	0.00
Target Compounds						
10) Phenol	6.52	94	32647	2.054	ng	88
49) Dimethylphthalate	9.63	163	105974	3.919	ng	99
70) Phenanthrene	11.43	178	63387	2.046	ng	100
74) Fluoranthene	12.62	202	107570	3.450	ng	96
77) Pyrene	12.85	202	97100	3.504	ng	98
80) Benzo(a)anthracene	14.04	228	45738	2.058	ng	97
82) Chrysene	14.07	228	48090	2.147	ng	100
87) Benzo(b)fluoranthene	15.10	252	59797m	2.568	ng	
89) Benzo(a)pyrene	15.47	252	42693	2.037	ng	# 95

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

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