

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072318\
 Data File : BF107611.D
 Acq On : 24 Jul 2018 9:38
 Operator : JU/SJ
 Sample : J4030-08DL 20X
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB7-12-7DL

Manual Integrations
 APPROVED

Sohil
 7/24/2018 5:26:28 PM

Quant Time: Jul 24 10:54:23 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF072018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 20 16:29:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.98	152	228675	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	900672	20.00	ng	0.00
38) Acenaphthene-d10	10.02	164	348354	20.00	ng	0.00
63) Phenanthrene-d10	11.52	188	639552	20.00	ng	0.00
75) Chrysene-d12	14.17	240	523562	20.00	ng	0.00
86) Perylene-d12	15.71	264	409383	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	18862m	1.33	ng	0.00
7) Phenol-d6	6.61	99	77228m	4.52	ng	0.00
23) Nitrobenzene-d5	7.54	82	44191m	3.31	ng	-0.01
41) 2,4,6-Tribromophenol	10.82	330	1659	0.42	ng	0.00
44) 2-Fluorobiphenyl	0.00	172	0d	0.00	ng	
78) Terphenyl-d14	13.10	244	87715	4.29	ng	0.00
Target Compounds						
70) Phenanthrene	11.54	178	76318	2.449	ng	99
74) Fluoranthene	12.74	202	126829	3.792	ng	96
77) Pyrene	12.97	202	115775	3.233	ng	98
83) Bis(2-ethylhexyl)phthalate	14.12	149	596491	27.449	ng	99
87) Benzo(b)fluoranthene	15.25	252	45736m	2.041	ng	

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

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