

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071718\
 Data File : BF107444.D
 Acq On : 17 Jul 2018 12:54
 Operator : JU/SJ
 Sample : J3819-05 2X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AU-05-071318-C

Manual Integrations
 APPROVED

Sohil
 7/18/2018 10:43:51 AM

Quant Time: Jul 17 14:26:25 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 17 11:29:08 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.98	152	154782	20.00	ng	0.00
21) Naphthalene-d8	8.27	136	584073	20.00	ng	0.00
38) Acenaphthene-d10	10.02	164	296490	20.00	ng	0.00
63) Phenanthrene-d10	11.52	188	560758	20.00	ng	0.00
75) Chrysene-d12	14.17	240	469733	20.00	ng	0.00
86) Perylene-d12	15.71	264	400965	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	458585	45.00	ng	0.00
7) Phenol-d6	6.60	99	592018	44.93	ng	0.00
23) Nitrobenzene-d5	7.54	82	425239	38.82	ng	-0.01
41) 2,4,6-Tribromophenol	10.82	330	130142	49.96	ng	0.00
44) 2-Fluorobiphenyl	9.34	172	702377	32.85	ng	-0.01
78) Terphenyl-d14	13.11	244	626843	33.16	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.73	163	128048	5.436	ng	96
53) 2,4-Dinitrophenol	10.06	184	585	2.792	ng	# 42
70) Phenanthrene	11.54	178	61438	2.198	ng	97
74) Fluoranthene	12.74	202	100119	3.152	ng	95
77) Pyrene	12.97	202	100313	3.140	ng	96
87) Benzo(b)fluoranthene	15.24	252	72252m	3.079	ng	
88) Benzo(k)fluoranthene	15.24	252	99569	4.347	ng	# 86
89) Benzo(a)pyrene	15.63	252	51128	2.374	ng	# 93

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

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