

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF072618\
 Data File : BF107717.D
 Acq On : 26 Jul 2018 23:10
 Operator : JU/SJ
 Sample : J4109-08
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-10(10-12)

Manual Integrations
 APPROVED

Sohil
 7/27/2018 2:54:13 PM

Quant Time: Jul 27 05:34:37 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	6.97	152	190028	20.00	ng	-0.01
21) Naphthalene-d8	8.24	136	784984	20.00	ng	-0.02
38) Acenaphthene-d10	10.00	164	348106	20.00	ng	-0.02
63) Phenanthrene-d10	11.50	188	615645	20.00	ng	-0.02
75) Chrysene-d12	14.14	240	432369	20.00	ng	-0.02
86) Perylene-d12	15.67	264	475537	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	1346057	113.89	ng	0.00
7) Phenol-d6	6.60	99	1619530	114.12	ng	-0.01
23) Nitrobenzene-d5	7.52	82	1059961	91.13	ng	-0.02
41) 2,4,6-Tribromophenol	10.80	330	487362	123.12	ng	-0.02
44) 2-Fluorobiphenyl	9.31	172	1879985	95.70	ng	-0.02
78) Terphenyl-d14	13.08	244	1356191	80.30	ng	-0.02
Target Compounds						
49) Dimethylphthalate	9.71	163	311082	11.857	ng	99
74) Fluoranthene	12.71	202	113875	3.537	ng	95
77) Pyrene	12.94	202	105203	3.557	ng	98
80) Benzo(a)anthracene	14.14	228	65424	2.582	ng	94
82) Chrysene	14.17	228	57982	2.382	ng	96
87) Benzo(b)fluoranthene	15.22	252	72141m	2.772	ng	
89) Benzo(a)pyrene	15.61	252	61806	2.596	ng	95

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

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