

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF072618\
 Data File : BF107709.D
 Acq On : 26 Jul 2018 19:32
 Operator : JU/SJ
 Sample : J4095-13
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CON-SB-104-05-06-07-0-11

Manual Integrations
 APPROVED

Sohil
 7/27/2018 2:53:53 PM

Quant Time: Jul 27 10:38:45 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	6.96	152	173321	20.00	ng	-0.02
21) Naphthalene-d8	8.24	136	699696	20.00	ng	-0.02
38) Acenaphthene-d10	10.00	164	330688	20.00	ng	-0.02
63) Phenanthrene-d10	11.49	188	654273	20.00	ng	-0.02
75) Chrysene-d12	14.14	240	398902	20.00	ng	-0.02
86) Perylene-d12	15.68	264	454878	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	844919	78.38	ng	-0.01
7) Phenol-d6	6.60	99	1044398	80.69	ng	-0.02
23) Nitrobenzene-d5	7.52	82	685437	66.11	ng	-0.02
41) 2,4,6-Tribromophenol	10.80	330	346929	92.26	ng	-0.02
44) 2-Fluorobiphenyl	9.31	172	1298215	62.31	ng	-0.02
78) Terphenyl-d14	13.08	244	1084795	69.62	ng	-0.02
Target Compounds						
49) Dimethylphthalate	9.71	163	121113	4.860	ng	99
70) Phenanthrene	11.52	178	106454	3.339	ng	99
74) Fluoranthene	12.71	202	168276	4.919	ng	95
77) Pyrene	12.94	202	144676	5.302	ng	97
80) Benzo(a)anthracene	14.14	228	60103	2.571	ng	94
82) Chrysene	14.17	228	56514	2.517	ng	98
85) Indeno(1,2,3-cd)pyrene	17.25	276	38979	2.048	ng	# 90
87) Benzo(b)fluoranthene	15.22	252	71922m	2.889	ng	
89) Benzo(a)pyrene	15.61	252	62380	2.739	ng	# 93
91) Benzo(g,h,i)perylene	17.72	276	40622	2.092	ng	# 90

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

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