

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072718\
 Data File : BF107758.D
 Acq On : 27 Jul 2018 20:49
 Operator : JU/SJ
 Sample : J4121-06
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GW-03(8-10)

Manual Integrations
 APPROVED

Sohil
 7/30/2018 2:03:04 PM

Quant Time: Jul 28 02:13:06 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	167469	20.00	ng	-0.02
21) Naphthalene-d8	8.24	136	663319	20.00	ng	-0.02
38) Acenaphthene-d10	10.00	164	269883	20.00	ng	-0.02
63) Phenanthrene-d10	11.49	188	471966	20.00	ng	-0.02
75) Chrysene-d12	14.15	240	432724	20.00	ng	-0.02
86) Perylene-d12	15.68	264	400907	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	988508	94.90	ng	-0.01
7) Phenol-d6	6.60	99	1282446	102.54	ng	-0.02
23) Nitrobenzene-d5	7.52	82	779267	79.28	ng	-0.02
41) 2,4,6-Tribromophenol	10.80	330	327753	106.80	ng	-0.02
44) 2-Fluorobiphenyl	9.31	172	1227039	75.49	ng	-0.02
78) Terphenyl-d14	13.08	244	939391	55.58	ng	-0.02
Target Compounds						
49) Dimethylphthalate	9.71	163	95125	4.677	ng	99
70) Phenanthrene	11.52	178	100960	4.390	ng	98
71) Anthracene	11.57	178	48854	2.110	ng	97
74) Fluoranthene	12.72	202	241818	9.798	ng	96
77) Pyrene	12.94	202	204921	6.923	ng	98
80) Benzo(a)anthracene	14.14	228	129456	5.105	ng	96
82) Chrysene	14.17	228	114179	4.687	ng	97
87) Benzo(b)fluoranthene	15.22	252	111040m	5.061	ng	
88) Benzo(k)fluoranthene	15.24	252	66878m	3.111	ng	
89) Benzo(a)pyrene	15.61	252	93869	4.677	ng	97
91) Benzo(g,h,i)perylene	17.73	276	34609	2.022	ng	# 93

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

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