

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072018\
 Data File : BF107547.D
 Acq On : 21 Jul 2018 1:49
 Operator : JU/SJ
 Sample : J4062-01 2X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-3

Manual Integrations
 APPROVED

Sohil
 7/23/2018 12:44:46 PM

Quant Time: Jul 21 05:09:17 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	325247	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	1240303	20.00	ng	0.00
38) Acenaphthene-d10	10.02	164	476401	20.00	ng	0.00
63) Phenanthrene-d10	11.51	188	832132	20.00	ng	0.00
75) Chrysene-d12	14.16	240	752952	20.00	ng	0.00
86) Perylene-d12	15.69	264	598416	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	924489	45.70	ng	0.00
7) Phenol-d6	6.60	99	1164113	47.93	ng	-0.01
23) Nitrobenzene-d5	7.54	82	654915	35.64	ng	-0.01
41) 2,4,6-Tribromophenol	10.81	330	234042	43.20	ng	-0.01
44) 2-Fluorobiphenyl	9.33	172	1114035	32.53	ng	0.00
78) Terphenyl-d14	13.09	244	910451	30.96	ng	0.00
Target Compounds						
70) Phenanthrene	11.54	178	135154	3.333	ng	98
74) Fluoranthene	12.72	202	317384	7.294	ng	99
77) Pyrene	12.96	202	269024	5.223	ng	97
80) Benzo(a)anthracene	14.15	228	144840	3.283	ng	97
82) Chrysene	14.18	228	141221	3.332	ng	98
87) Benzo(b)fluoranthene	15.24	252	162829m	4.972	ng	
89) Benzo(a)pyrene	15.62	252	94962	3.170	ng	# 91
91) Benzo(g,h,i)perylene	17.74	276	56580	2.215	ng	# 92

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

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