

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070218\
 Data File : BF107038.D
 Acq On : 2 Jul 2018 12:23
 Operator : JU/SJ
 Sample : J3524-03DLRE 10X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AREA-3(D)DLRE

Manual Integrations
 APPROVED

Sohil
 7/3/2018 5:29:44 PM

Quant Time: Jul 02 14:27:56 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 22 11:34:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	75125	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	284687	20.00	ng	0.00
38) Acenaphthene-d10	10.00	164	129003	20.00	ng	0.00
63) Phenanthrene-d10	11.50	188	222481	20.00	ng	0.00
75) Chrysene-d12	14.15	240	229209	20.00	ng	0.00
86) Perylene-d12	15.70	264	174640	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	15744	3.55	ng	0.00
7) Phenol-d6	6.61	99	19449	3.51	ng	0.00
23) Nitrobenzene-d5	7.52	82	10990	2.15	ng	-0.01
41) 2,4,6-Tribromophenol	10.80	330	4183	2.80	ng	0.00
44) 2-Fluorobiphenyl	9.31	172	23693	-14.09	ng	-0.02
78) Terphenyl-d14	13.07	244	24551	2.59	ng	-0.01
Target Compounds						Qvalue
31) Naphthalene	8.26	128	351858	26.436	ng	99
37) 2-Methylnaphthalene	8.95	142	202798	22.987	ng	96
45) 1,1'-Biphenyl	9.41	154	21199	2.062	ng	100
48) Acenaphthylene	9.86	152	50754	3.972	ng	# 94
51) Acenaphthene	10.03	154	69537	8.643	ng	98
54) Dibenzofuran	10.21	168	39854	3.618	ng	93
57) Fluorene	10.55	166	116119	13.283	ng	99
70) Phenanthrene	11.52	178	465943	43.421	ng	99
71) Anthracene	11.57	178	125887	11.493	ng	99
74) Fluoranthene	12.72	202	307897	26.033	ng	94
77) Pyrene	12.95	202	396656	26.243	ng	99
80) Benzo(a)anthracene	14.14	228	182446	13.060	ng	93
82) Chrysene	14.18	228	150769	11.731	ng	98
85) Indeno(1,2,3-cd)pyrene	17.28	276	36949	3.388	ng	# 89
87) Benzo(b)fluoranthene	15.24	252	121105m	11.163	ng	
88) Benzo(k)fluoranthene	15.27	252	33752m	3.267	ng	
89) Benzo(a)pyrene	15.63	252	98850	10.250	ng	95
91) Benzo(g,h,i)perylene	17.77	276	39164	4.902	ng	# 87

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

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