

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050318\
 Data File : BF105086.D
 Acq On : 3 May 2018 21:25
 Operator : JU/SJ
 Sample : J2667-08
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 137-132-165

Manual Integrations
 APPROVED

Sohil
 5/4/2018 1:46:49 PM

Quant Time: May 04 09:51:52 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 02 17:25:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.78	152	106442	20.00	ng	0.00
21) Naphthalene-d8	8.06	136	451012	20.00	ng	-0.01
38) Acenaphthene-d10	9.82	164	212285	20.00	ng	-0.02
63) Phenanthrene-d10	11.31	188	355283	20.00	ng	-0.01
75) Chrysene-d12	14.01	240	211586	20.00	ng	-0.02
86) Perylene-d12	15.57	264	194154	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	759179	109.06	ng	0.00
7) Phenol-d6	6.43	99	920239	107.59	ng	0.00
23) Nitrobenzene-d5	7.35	82	560493	81.15	ng	-0.02
41) 2,4,6-Tribromophenol	10.62	330	243287	110.48	ng	-0.02
44) 2-Fluorobiphenyl	9.14	172	1067150	79.41	ng	-0.02
78) Terphenyl-d14	12.90	244	884327	95.29	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.54	163	87740	5.241	ng	99
70) Phenanthrene	11.33	178	182099	9.204	ng	99
74) Fluoranthene	12.53	202	233740	11.307	ng	97
77) Pyrene	12.76	202	189770	12.100	ng	99
80) Benzo(a)anthracene	14.00	228	91785	7.261	ng	98
82) Chrysene	14.03	228	83485	6.430	ng	98
85) Indeno(1,2,3-cd)pyrene	17.05	276	43546	3.948	ng	# 91
87) Benzo(b)fluoranthene	15.13	252	102985m	8.398	ng	
88) Benzo(k)fluoranthene	15.16	252	42445m	3.666	ng	
89) Benzo(a)pyrene	15.50	252	79689	7.263	ng	97
91) Benzo(g,h,i)perylene	17.50	276	42833	4.906	ng	97

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

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