

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050218\
 Data File : BF105024.D
 Acq On : 2 May 2018 12:17
 Operator : JU/SJ
 Sample : J2157-21
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-02-043018-E

Manual Integrations
 APPROVED

Sohil
 5/3/2018 5:00:44 PM

Quant Time: May 03 10:04:41 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	6.79	152	98687	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	408295	20.00	ng	0.00
38) Acenaphthene-d10	9.83	164	183131	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	310827	20.00	ng	0.00
75) Chrysene-d12	14.02	240	218478	20.00	ng	-0.01
86) Perylene-d12	15.59	264	206997	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	626229	97.03	ng	0.00
7) Phenol-d6	6.43	99	759779	95.81	ng	0.00
23) Nitrobenzene-d5	7.36	82	457673	73.20	ng	-0.01
41) 2,4,6-Tribromophenol	10.63	330	196800	103.60	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	859369	74.13	ng	0.00
78) Terphenyl-d14	12.92	244	680816	71.04	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.55	163	97192	6.730	ng	100
70) Phenanthrene	11.35	178	74378	4.297	ng	97
74) Fluoranthene	12.54	202	107707	5.955	ng	96
77) Pyrene	12.77	202	106341	6.567	ng	98
80) Benzo(a)anthracene	14.01	228	49652	3.804	ng	96
82) Chrysene	14.05	228	53264	3.973	ng	95
85) Indeno(1,2,3-cd)pyrene	17.08	276	26443	2.322	ng	# 90
87) Benzo(b)fluoranthene	15.16	252	68311m	5.225	ng	
89) Benzo(a)pyrene	15.53	252	46666m	3.989	ng	
91) Benzo(g,h,i)perylene	17.53	276	26628	2.861	ng	94

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

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