

Data Path : Z:\HPCHEM1\BNA F\DATA\BF041218\
 Data File : BF104450.D
 Acq On : 12 Apr 2018 14:55
 Operator : JU/SJ
 Sample : J2165-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-03-041118

Manual Integrations
 APPROVED

Sohil
 4/13/2018 4:43:57 PM

Quant Time: Apr 13 10:03:51 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 12 17:07:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	145719	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	577917	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	248018	20.00	ng	0.00
63) Phenanthrene-d10	11.33	188	372981	20.00	ng	0.00
75) Chrysene-d12	13.98	240	280965	20.00	ng	-0.01
86) Perylene-d12	15.45	264	227859	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	931640	97.76	ng	0.01
7) Phenol-d6	6.45	99	1119719	95.63	ng	0.00
23) Nitrobenzene-d5	7.37	82	719120	81.25	ng	0.00
41) 2,4,6-Tribromophenol	10.64	330	234499	91.15	ng	0.00
44) 2-Fluorobiphenyl	9.17	172	1105341	70.40	ng	0.00
78) Terphenyl-d14	12.93	244	799300	64.86	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.56	163	78634	4.021	ng	99
70) Phenanthrene	11.36	178	64628	3.111	ng	98
74) Fluoranthene	12.55	202	112569	5.187	ng	97
77) Pyrene	12.78	202	107835	5.178	ng	98
80) Benzo(a)anthracene	13.97	228	55143	3.285	ng	97
82) Chrysene	14.00	228	57380	3.328	ng	96
87) Benzo(b)fluoranthene	15.02	252	71442m	4.964	ng	
89) Benzo(a)pyrene	15.39	252	42013	3.263	ng	# 86
91) Benzo(g,h,i)perylene	17.34	276	24391	2.381	ng	# 93

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

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