

Data Path : Z:\HPCHEM1\BNA F\DATA\BF031218\
 Data File : BF103574.D
 Acq On : 12 Mar 2018 11:46
 Operator : SJ/JU
 Sample : J1776-08
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-1-6FT

Manual Integrations
 APPROVED

Sohil
 3/13/2018 4:12:45 PM

Quant Time: Mar 12 16:28:35 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF022218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 09 11:47:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	141927	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	599819	20.00	ng	0.00
38) Acenaphthene-d10	9.90	164	252020	20.00	ng	-0.01
63) Phenanthrene-d10	11.40	188	401509	20.00	ng	0.00
75) Chrysene-d12	14.06	240	225495	20.00	ng	0.00
86) Perylene-d12	15.57	264	212847	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	762311	81.23	ng	0.01
7) Phenol-d6	6.50	99	982395	85.55	ng	0.00
23) Nitrobenzene-d5	7.43	82	612875	58.35	ng	-0.01
41) 2,4,6-Tribromophenol	10.70	330	134276	62.95	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	862731	53.86	ng	0.00
78) Terphenyl-d14	12.98	244	575991	61.40	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.62	163	71981	3.560	ng	98
70) Phenanthrene	11.42	178	132352	5.990	ng	99
74) Fluoranthene	12.62	202	204514	9.211	ng	97
77) Pyrene	12.85	202	179277	10.197	ng	99
80) Benzo(a)anthracene	14.04	228	65499m	4.477	ng	
82) Chrysene	14.09	228	66650	4.692	ng	96
85) Indeno(1,2,3-cd)pyrene	17.12	276	40334	3.586	ng	# 92
87) Benzo(b)fluoranthene	15.13	252	90221m	6.447	ng	
88) Benzo(k)fluoranthene	15.15	252	29537m	2.241	ng	
89) Benzo(a)pyrene	15.51	252	65371	5.231	ng	# 86
91) Benzo(g,h,i)perylene	17.60	276	40727	4.315	ng	# 84

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

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