

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121217\
 Data File : BF101351.D
 Acq On : 13 Dec 2017 00:55
 Operator : SJ/JU
 Sample : I6822-12
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-8(13-15)

Manual Integrations
 APPROVED

Sohil
 12/13/2017 12:08:29 PM

Quant Time: Dec 13 05:34:46 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 08 18:17:31 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	47501	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	179738	20.00	ng	0.00
38) Acenaphthene-d10	9.86	164	76491	20.00	ng	0.00
63) Phenanthrene-d10	11.36	188	128190	20.00	ng	0.00
75) Chrysene-d12	14.00	240	107686	20.00	ng	0.00
86) Perylene-d12	15.48	264	89495	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	324381	112.84	ng	0.00
7) Phenol-d6	6.46	99	386676	110.79	ng	0.00
23) Nitrobenzene-d5	7.39	82	220876	73.31	ng	0.00
41) 2,4,6-Tribromophenol	10.66	330	83898	91.31	ng	0.00
44) 2-Fluorobiphenyl	9.18	172	430434	76.99	ng	0.00
78) Terphenyl-d14	12.94	244	345232	70.12	ng	0.00
Target Compounds						
10) Phenol	6.47	94	13598	3.504	ng	87
49) Dimethylphthalate	9.57	163	71852	11.648	ng	# 99
70) Phenanthrene	11.38	178	42085	5.962	ng	96
71) Anthracene	11.43	178	14447	2.022	ng	93
74) Fluoranthene	12.57	202	55638	7.110	ng	93
77) Pyrene	12.80	202	55590	7.481	ng	99
80) Benzo(a)anthracene	13.99	228	39392	5.794	ng	96
82) Chrysene	14.03	228	35035	5.548	ng	96
85) Indeno(1,2,3-cd)pyrene	16.97	276	31222	5.562	ng	# 88
87) Benzo(b)fluoranthene	15.05	252	53505m	9.981	ng	
88) Benzo(k)fluoranthene	15.07	252	16861m	3.193	ng	
89) Benzo(a)pyrene	15.42	252	46136	9.467	ng	96
90) Dibenzo(a,h)anthracene	16.97	278	8941	2.081	ng	92
91) Benzo(a,h,i)perylene	17.42	276	27887	6.888	ng	# 89

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

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