

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121217\
 Data File : BF101348.D
 Acq On : 12 Dec 2017 23:35
 Operator : SJ/JU
 Sample : I6814-01
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PHMHC-WC1

Manual Integrations
 APPROVED

Sohil
 12/13/2017 12:08:09 PM

Quant Time: Dec 13 04:41:40 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	40246	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	154466	20.00	ng	0.00
38) Acenaphthene-d10	9.86	164	66722	20.00	ng	0.00
63) Phenanthrene-d10	11.35	188	115803	20.00	ng	0.00
75) Chrysene-d12	14.00	240	94101	20.00	ng	0.00
86) Perylene-d12	15.49	264	78376	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	252708	103.76	ng	0.00
7) Phenol-d6	6.46	99	307654	104.04	ng	0.00
23) Nitrobenzene-d5	7.39	82	181665	70.16	ng	-0.01
41) 2,4,6-Tribromophenol	10.66	330	68948	86.02	ng	0.00
44) 2-Fluorobiphenyl	9.18	172	333533	68.39	ng	0.00
78) Terphenyl-d14	12.94	244	258911	60.18	ng	0.00
Target Compounds						
10) Phenol	6.47	94	10218	3.108	ng	87
48) Acenaphthylene	9.72	152	20607	2.888	ng	98
49) Dimethylphthalate	9.57	163	24855	4.619	ng	# 96
70) Phenanthrene	11.38	178	45674	7.162	ng	97
71) Anthracene	11.43	178	16542	2.563	ng	99
74) Fluoranthene	12.57	202	75268	10.648	ng	94
77) Pyrene	12.80	202	83697	12.890	ng	97
80) Benzo(a)anthracene	14.00	228	47362	7.972	ng	94
82) Chrysene	14.03	228	42830	7.762	ng	98
85) Indeno(1,2,3-cd)pyrene	16.97	276	28464	5.802	ng	# 90
87) Benzo(b)fluoranthene	15.05	252	68845m	14.665	ng	
88) Benzo(k)fluoranthene	15.08	252	13906m	3.007	ng	
89) Benzo(a)pyrene	15.42	252	45816	10.735	ng	96
90) Dibenzo(a,h)anthracene	16.97	278	10084	2.680	ng	# 66
91) Benzo(a,h,i)perylene	17.43	276	34854	9.829	ng	94

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

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