

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122017\
 Data File : BF101582.D
 Acq On : 20 Dec 2017 21:47
 Operator : SJ/JU
 Sample : I6981-07 2X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PHMHB-WCS

Manual Integrations
 APPROVED

Sohil
 12/21/2017 12:07:06 PM

Quant Time: Dec 21 04:52:03 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 20 14:25:04 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	142602	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	521409	20.00	ng	0.00
38) Acenaphthene-d10	9.83	164	193603	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	300644	20.00	ng	0.00
75) Chrysene-d12	13.97	240	303934	20.00	ng	0.00
86) Perylene-d12	15.44	264	293770	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	533576	55.74	ng	0.00
7) Phenol-d6	6.43	99	637067	53.47	ng	0.00
23) Nitrobenzene-d5	7.36	82	374537	39.99	ng	0.00
41) 2,4,6-Tribromophenol	10.63	330	84760	45.44	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	522525	41.87	ng	0.00
78) Terphenyl-d14	12.90	244	395140	31.34	ng	-0.01
Target Compounds						
10) Phenol	6.45	94	29127	2.145	ng	Qvalue # 69
51) Acenaphthene	9.86	154	50012	4.012	ng	98
54) Dibenzofuran	10.04	168	37628	2.375	ng	98
57) Fluorene	10.38	166	67837	5.062	ng	99
70) Phenanthrene	11.35	178	288596	17.261	ng	98
71) Anthracene	11.40	178	118532	6.941	ng	97
74) Fluoranthene	12.54	202	403591	22.998	ng	98
77) Pyrene	12.77	202	387707	16.997	ng	100
80) Benzo(a)anthracene	13.96	228	189176	9.426	ng	97
82) Chrysene	14.00	228	180453	9.523	ng	97
85) Indeno(1,2,3-cd)pyrene	16.93	276	77875	4.615	ng	# 91
87) Benzo(b)fluoranthene	15.02	252	197792m	11.046	ng	
88) Benzo(k)fluoranthene	15.04	252	58456m	3.358	ng	
89) Benzo(a)pyrene	15.38	252	157334	9.532	ng	# 91
91) Benzo(g,h,i)perylene	17.37	276	80568	5.741	ng	# 94

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

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