

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122717\
 Data File : BF101783.D
 Acq On : 28 Dec 2017 00:35
 Operator : SJ/JU
 Sample : I7064-13
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SAMPLE-7

Manual Integrations
 APPROVED

Sohil
 12/28/2017 2:03:16 PM

Quant Time: Dec 28 06:04:43 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 26 10:48:01 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	163901	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	622131	20.00	ng	0.00
38) Acenaphthene-d10	9.83	164	247533	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	384739	20.00	ng	0.00
75) Chrysene-d12	13.97	240	304851	20.00	ng	0.00
86) Perylene-d12	15.44	264	263009	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	1130874	102.78	ng	0.01
7) Phenol-d6	6.44	99	1326384	96.86	ng	0.00
23) Nitrobenzene-d5	7.36	82	815449	72.96	ng	0.00
41) 2,4,6-Tribromophenol	10.63	330	234361	98.26	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	1124296	70.46	ng	0.00
78) Terphenyl-d14	12.90	244	748701	59.21	ng	0.00
Target Compounds						
10) Phenol	6.45	94	71494	4.581	ng	73
49) Dimethylphthalate	9.55	163	134675	6.764	ng	98
54) Dibenzofuran	10.04	168	42577	2.102	ng	98
57) Fluorene	10.38	166	39398	2.299	ng	96
70) Phenanthrene	11.35	178	615840	28.783	ng	99
71) Anthracene	11.40	178	124903	5.715	ng	96
74) Fluoranthene	12.54	202	641187	28.550	ng	98
77) Pyrene	12.77	202	565297	24.708	ng	99
80) Benzo(a)anthracene	13.96	228	274166	13.620	ng	98
82) Chrysene	14.00	228	239904	12.622	ng	99
85) Indeno(1,2,3-cd)pyrene	16.94	276	73547	4.345	ng	89
87) Benzo(b)fluoranthene	15.02	252	210842m	13.152	ng	
88) Benzo(k)fluoranthene	15.04	252	89536m	5.744	ng	
89) Benzo(a)pyrene	15.39	252	154261	10.438	ng	96
91) Benzo(a,h,i)perylene	17.40	276	61159	4.868	ng	95

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

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