

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122717\
 Data File : BF101778.D
 Acq On : 27 Dec 2017 22:22
 Operator : SJ/JU
 Sample : I7064-03 2X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SAMPLE-2

Manual Integrations
 APPROVED

Sohil
 12/28/2017 2:03:02 PM

Quant Time: Dec 28 05:44:38 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	162470	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	624160	20.00	ng	0.00
38) Acenaphthene-d10	9.83	164	255293	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	412245	20.00	ng	0.00
75) Chrysene-d12	13.97	240	299290	20.00	ng	0.00
86) Perylene-d12	15.44	264	277952	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	661480	60.65	ng	0.00
7) Phenol-d6	6.43	99	794662	58.54	ng	0.00
23) Nitrobenzene-d5	7.36	82	475275	42.39	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	143129	58.18	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	806391	49.00	ng	0.00
78) Terphenyl-d14	12.90	244	554401	44.66	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.55	163	55402	2.698	ng	99
70) Phenanthrene	11.34	178	455767	19.880	ng	99
71) Anthracene	11.40	178	121376	5.183	ng	99
72) Carbazole	11.57	167	53757	2.452	ng	98
74) Fluoranthene	12.54	202	618397	25.698	ng	97
77) Pyrene	12.77	202	541982	24.129	ng	99
80) Benzo(a)anthracene	13.96	228	285885	14.466	ng	99
82) Chrysene	14.00	228	274746m	14.724	ng	
85) Indeno(1,2,3-cd)pyrene	16.94	276	114658	6.900	ng	# 89
87) Benzo(b)fluoranthene	15.02	252	309823m	18.287	ng	
88) Benzo(k)fluoranthene	15.04	252	115506m	7.012	ng	
89) Benzo(a)pyrene	15.39	252	207029	13.256	ng	98
90) Dibenzo(a,h)anthracene	16.93	278	27399	2.043	ng	# 88
91) Benzo(g,h,i)perylene	17.39	276	105508	7.946	ng	96

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

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