

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050817\
 Data File : BF094988.D
 Acq On : 8 May 2017 19:32
 Operator : SJ/MA
 Sample : I2947-14
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-05-0-2.0

Manual Integrations
 APPROVED

mohammad
 5/9/2017 5:02:34 PM

Quant Time: May 09 02:08:21 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 03 17:24:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	94096	20.00	ng	-0.02
21) Naphthalene-d8	9.74	136	393048	20.00	ng	-0.02
38) Acenaphthene-d10	12.56	164	197131	20.00	ng	-0.02
63) Phenanthrene-d10	14.97	188	305554	20.00	ng	-0.02
75) Chrysene-d12	18.63	240	243540	20.00	ng	-0.02
86) Perylene-d12	20.30	264	213407	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.65	112	667652	118.30	ng	-0.02
7) Phenol-d6	7.25	99	748523	110.53	ng	-0.03
23) Nitrobenzene-d5	8.61	82	444637	71.34	ng	-0.03
41) 2,4,6-Tribromophenol	13.86	330	162604	100.72	ng	-0.02
44) 2-Fluorobiphenyl	11.51	172	825598	60.28	ng	-0.02
78) Terphenyl-d14	17.29	244	571107	51.65	ng	-0.02
Target Compounds						Qvalue
10) Phenol	7.27	94	15861	2.22	ng	82
49) Dimethylphthalate	12.20	163	164902	10.19	ng	98
74) Fluoranthene	16.75	202	88839	4.93	ng	99
77) Pyrene	17.05	202	91955	5.05	ng	99
80) Benzo(a)anthracene	18.62	228	49468	3.49	ng	92
82) Chrysene	18.67	228	65912	4.81	ng	98
85) Indeno(1,2,3-cd)pyrene	21.60	276	35844	3.22	ng	98
87) Benzo(b)fluoranthene	19.89	252	83645m	6.34	ng	
88) Benzo(k)fluoranthene	19.92	252	31733m	2.49	ng	
89) Benzo(a)pyrene	20.25	252	59730	4.91	ng	99
91) Benzo(g,h,i)perylene	21.98	276	38373	3.89	ng	95

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

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