

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070717\
 Data File : BF096573.D
 Acq On : 7 Jul 2017 21:48
 Operator : SJ/MA
 Sample : I4063-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NB-01-070517-C

Manual Integrations
 APPROVED

mohammad
 7/10/2017 2:22:36 PM

Quant Time: Jul 08 00:49:42 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 07 13:02:33 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	113641	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	506989	20.00	ng	-0.01
38) Acenaphthene-d10	9.73	164	219342	20.00	ng	0.00
63) Phenanthrene-d10	11.20	188	348204	20.00	ng	0.00
75) Chrysene-d12	13.83	240	237865	20.00	ng	-0.01
86) Perylene-d12	15.23	264	238709	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.31	112	732908	95.19	ng	0.00
7) Phenol-d6	6.33	99	914994	96.67	ng	-0.01
23) Nitrobenzene-d5	7.26	82	563259	66.76	ng	-0.01
41) 2,4,6-Tribromophenol	10.52	330	177782	103.76	ng	0.00
44) 2-Fluorobiphenyl	9.06	172	971197	75.70	ng	0.00
78) Terphenyl-d14	12.79	244	662769	59.54	ng	0.00
Target Compounds						
10) Phenol	6.35	94	32390	3.004	ng	94
49) Dimethylphthalate	9.45	163	133321	8.142	ng	98
70) Phenanthrene	11.23	178	50430	2.679	ng	97
74) Fluoranthene	12.41	202	100620	5.452	ng	99
77) Pyrene	12.63	202	98894	5.180	ng	98
80) Benzo(a)anthracene	13.82	228	40161	2.916	ng	97
82) Chrysene	13.86	228	45956	3.186	ng	97
85) Indeno(1,2,3-cd)pyrene	16.57	276	24359	2.139	ng	96
87) Benzo(b)fluoranthene	14.84	252	56051m	3.747	ng	
89) Benzo(a)pyrene	15.17	252	42399	3.175	ng	# 95
91) Benzo(g,h,i)perylene	16.97	276	26017	2.390	ng	98

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

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