

Data Path : Z:\HPCHEM1\BNA F\DATA\BF041317\
 Data File : BF094403.D
 Acq On : 13 Apr 2017 21:07
 Operator : SJ/MA
 Sample : I2662-01
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AREA-1-COMP

Manual Integrations
 APPROVED

mohammad
 4/14/2017 1:37:01 PM

Quant Time: Apr 14 07:12:35 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF032217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 14 01:21:26 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.79	152	162962	20.00	ng	-0.04
21) Naphthalene-d8	7.29	136	625933	20.00	ng	-0.05
38) Acenaphthene-d10	9.42	164	266718	20.00	ng	-0.05
63) Phenanthrene-d10	11.24	188	410827	20.00	ng	-0.05
75) Chrysene-d12	14.49	240	302719	20.00	ng	-0.06
86) Perylene-d12	16.12	264	252304	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	4.33	112	561832	58.23	ng	-0.05
7) Phenol-d6	5.42	99	903602	75.70	ng	-0.05
23) Nitrobenzene-d5	6.43	82	483758	44.11	ng	-0.05
41) 2,4,6-Tribromophenol	10.40	330	205607	97.55	ng	-0.05
44) 2-Fluorobiphenyl	8.61	172	1062268	60.75	ng	-0.05
78) Terphenyl-d14	13.22	244	770590	57.06	ng	-0.05
Target Compounds						Qvalue
10) Phenol	5.43	94	51331	3.78	ng	91
49) Dimethylphthalate	9.10	163	136310	7.19	ng	98
70) Phenanthrene	11.27	178	235705	10.31	ng	99
71) Anthracene	11.33	178	57288	2.43	ng	98
74) Fluoranthene	12.73	202	499161	21.33	ng	99
77) Pyrene	13.01	202	477250	21.72	ng	99
80) Benzo(a)anthracene	14.48	228	202149	11.45	ng	99
82) Chrysene	14.53	228	195897m	11.96	ng	
85) Indeno(1,2,3-cd)pyrene	17.41	276	101685	7.17	ng	95
87) Benzo(b)fluoranthene	15.72	252	230786m	14.92	ng	
88) Benzo(k)fluoranthene	15.74	252	90717m	6.00	ng	
89) Benzo(a)pyrene	16.06	252	174316	12.24	ng	99
91) Benzo(a,h,i)perylene	17.79	276	126071	10.37	ng	98

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

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