

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041317\
 Data File : BF094423.D
 Acq On : 14 Apr 2017 10:32
 Operator : SJ/MA
 Sample : I2672-01 5X
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-Z8-NSW

Manual Integrations
 APPROVED

mohammad
 4/14/2017 1:36:23 PM

Quant Time: Apr 14 10:56:35 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	137108	20.00	ng	-0.04
21) Naphthalene-d8	7.29	136	530246	20.00	ng	-0.05
38) Acenaphthene-d10	9.42	164	222693	20.00	ng	-0.05
63) Phenanthrene-d10	11.25	188	340871	20.00	ng	-0.05
75) Chrysene-d12	14.49	240	262332	20.00	ng	-0.06
86) Perylene-d12	16.12	264	200483	20.00	ng	-0.06
System Monitoring Compounds						
5) 2-Fluorophenol	4.33	112	200030	24.64	ng	-0.05
7) Phenol-d6	5.40	99	242267	24.12	ng	-0.06
23) Nitrobenzene-d5	6.43	82	139361	15.00	ng	-0.06
41) 2,4,6-Tribromophenol	10.39	330	28729	16.33	ng	-0.05
44) 2-Fluorobiphenyl	8.60	172	262844	18.00	ng	-0.05
78) Terphenyl-d14	13.22	244	163607	13.98	ng	-0.06
Target Compounds						Qvalue
49) Dimethylphthalate	9.10	163	39195	2.48	ng	98
70) Phenanthrene	11.27	178	187110	9.87	ng	98
71) Anthracene	11.33	178	43523	2.23	ng	97
74) Fluoranthene	12.73	202	144364	7.44	ng	99
77) Pyrene	13.01	202	170611	8.96	ng	98
80) Benzo(a)anthracene	14.48	228	59268	3.87	ng	99
82) Chrysene	14.52	228	57222	4.03	ng	97
87) Benzo(b)fluoranthene	15.72	252	43826m	3.57	ng	
89) Benzo(a)pyrene	16.06	252	35325	3.12	ng	99

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

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