

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041317\
 Data File : BF094424.D
 Acq On : 14 Apr 2017 11:00
 Operator : SJ/MA
 Sample : I2672-07 5X
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-AA10-SSW-2

Manual Integrations
 APPROVED

mohammad
 4/14/2017 1:36:17 PM

Quant Time: Apr 14 11:24:36 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	134111	20.00	ng	-0.04
21) Naphthalene-d8	7.29	136	517606	20.00	ng	-0.05
38) Acenaphthene-d10	9.42	164	216671	20.00	ng	-0.05
63) Phenanthrene-d10	11.25	188	328629	20.00	ng	-0.05
75) Chrysene-d12	14.49	240	255550	20.00	ng	-0.06
86) Perylene-d12	16.12	264	196062	20.00	ng	-0.06
System Monitoring Compounds						
5) 2-Fluorophenol	4.33	112	187447	23.61	ng	-0.05
7) Phenol-d6	5.40	99	232073	23.63	ng	-0.06
23) Nitrobenzene-d5	6.43	82	133747	14.75	ng	-0.06
41) 2,4,6-Tribromophenol	10.39	330	31647	18.48	ng	-0.05
44) 2-Fluorobiphenyl	8.60	172	251500	17.70	ng	-0.05
78) Terphenyl-d14	13.22	244	149943	13.15	ng	-0.06
Target Compounds						Qvalue
49) Dimethylphthalate	9.10	163	38135	2.48	ng	# 99
70) Phenanthrene	11.27	178	85440	4.67	ng	98
74) Fluoranthene	12.73	202	127977	6.84	ng	99
77) Pyrene	13.01	202	108382	5.84	ng	97
80) Benzo(a)anthracene	14.48	228	51410	3.45	ng	97
82) Chrysene	14.52	228	46900	3.39	ng	96
87) Benzo(b)fluoranthene	15.72	252	50924m	4.24	ng	
89) Benzo(a)pyrene	16.06	252	33963	3.07	ng	99

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

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