

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021017\
 Data File : BF092900.D
 Acq On : 10 Feb 2017 23:07
 Operator : SJ/MA
 Sample : I1777-04
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 38-SB-03-0-2

Manual Integrations
 APPROVED

mohammad
 2/13/2017 12:41:07 PM

Quant Time: Feb 11 00:33:33 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF013017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 10 23:31:14 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	159559	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	624457	20.00	ng	-0.01
38) Acenaphthene-d10	9.94	164	375137	20.00	ng	0.00
63) Phenanthrene-d10	11.42	188	635287	20.00	ng	0.00
75) Chrysene-d12	14.05	240	478548	20.00	ng	-0.01
86) Perylene-d12	15.51	264	371679	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	818570	88.01	ng	0.00
7) Phenol-d6	6.53	99	1134277	94.79	ng	0.00
23) Nitrobenzene-d5	7.46	82	731142	65.20	ng	-0.01
41) 2,4,6-Tribromophenol	10.73	330	215414	79.39	ng	-0.01
44) 2-Fluorobiphenyl	9.25	172	1409539	68.07	ng	-0.01
78) Terphenyl-d14	13.01	244	1343069	65.94	ng	0.00
Target Compounds						
10) Phenol	6.54	94	71363	5.10	ng	Qvalue # 72
49) Dimethylphthalate	9.65	163	89006	3.52	ng	99
70) Phenanthrene	11.45	178	86056	2.36	ng	97
74) Fluoranthene	12.64	202	157330	4.19	ng	93
77) Pyrene	12.87	202	161126	4.50	ng	97
80) Benzo(a)anthracene	14.04	228	75507	2.58	ng	98
82) Chrysene	14.08	228	58644m	2.14	ng	
87) Benzo(b)fluoranthene	15.08	252	73254m	2.99	ng	
89) Benzo(a)pyrene	15.44	252	46936	2.22	ng	# 93

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

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