

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062917\
 Data File : BF096311.D
 Acq On : 30 Jun 2017 00:15
 Operator : SJ/MA
 Sample : I3882-10 5X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 G47-1.5-3

Manual Integrations
 APPROVED

mohammad
 6/30/2017 2:52:17 PM

Quant Time: Jun 30 03:52:30 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF062717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 28 15:41:49 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	169535	20.00	ng	-0.01
21) Naphthalene-d8	8.05	136	643480	20.00	ng	0.00
38) Acenaphthene-d10	9.80	164	224514	20.00	ng	0.00
63) Phenanthrene-d10	11.29	188	341439	20.00	ng	0.00
75) Chrysene-d12	13.92	240	329128	20.00	ng	0.00
86) Perylene-d12	15.34	264	225458	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	248232	21.09	ng	0.00
7) Phenol-d6	6.39	99	298168	20.93	ng	0.00
23) Nitrobenzene-d5	7.32	82	151600	16.31	ng	-0.04
41) 2,4,6-Tribromophenol	10.59	330	41608	23.55	ng	0.00
44) 2-Fluorobiphenyl	9.13	172	260345	21.24	ng	-0.02
78) Terphenyl-d14	12.87	244	189234	14.35	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.52	163	33931	2.11	ng	# 97
74) Fluoranthene	12.49	202	118496	6.25	ng	96
77) Pyrene	12.72	202	115560	4.27	ng	99
80) Benzo(a)anthracene	13.91	228	49018	2.44	ng	98
82) Chrysene	13.94	228	55111	2.66	ng	94
87) Benzo(b)fluoranthene	14.93	252	57045m	3.85	ng	
89) Benzo(a)pyrene	15.28	252	28439	2.17	ng	# 71

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

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