

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070117\
 Data File : BF096388.D
 Acq On : 1 Jul 2017 19:36
 Operator : SJ/MA
 Sample : I3908-11
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 G-61-0-1.5

Manual Integrations
 APPROVED

mohammad
 7/3/2017 4:02:53 PM

Quant Time: Jul 03 02:38:58 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF070117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 03 01:43:12 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	153697	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	627147	20.00	ng	0.00
38) Acenaphthene-d10	9.76	164	222803	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	315318	20.00	ng	-0.01
75) Chrysene-d12	13.87	240	291707	20.00	ng	-0.01
86) Perylene-d12	15.28	264	263091	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	929306	89.24	ng	0.00
7) Phenol-d6	6.37	99	1140240	89.07	ng	0.00
23) Nitrobenzene-d5	7.29	82	691003	66.21	ng	-0.01
41) 2,4,6-Tribromophenol	10.55	330	148883	85.55	ng	0.00
44) 2-Fluorobiphenyl	9.09	172	1011553	77.63	ng	0.00
78) Terphenyl-d14	12.83	244	593793	43.50	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.48	163	278110	16.721	ng	99
70) Phenanthrene	11.26	178	43814	2.570	ng	98
74) Fluoranthene	12.45	202	63059	3.773	ng	99
77) Pyrene	12.67	202	58209	2.486	ng	99
80) Benzo(a)anthracene	13.86	228	36631	2.168	ng	91
82) Chrysene	13.89	228	42701	2.414	ng	# 95
87) Benzo(b)fluoranthene	14.88	252	53793m	3.263	ng	

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

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