

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060517\
 Data File : BF095692.D
 Acq On : 6 Jun 2017 1:22
 Operator : SJ/MA
 Sample : I3466-03
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-RO-4672-060217

Manual Integrations
 APPROVED

mohammad
 6/6/2017 5:13:04 PM

Quant Time: Jun 06 03:05:59 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 05 16:15:31 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	92528	20.00	ng	0.00
21) Naphthalene-d8	9.43	136	355884	20.00	ng	0.00
38) Acenaphthene-d10	12.25	164	148649	20.00	ng	0.00
63) Phenanthrene-d10	14.64	188	229549	20.00	ng	0.00
75) Chrysene-d12	18.35	240	205346	20.00	ng	0.00
86) Perylene-d12	20.02	264	154026	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	612658	105.51	ng	0.00
7) Phenol-d6	6.95	99	714119	102.73	ng	0.00
23) Nitrobenzene-d5	8.31	82	448533	78.27	ng	0.00
41) 2,4,6-Tribromophenol	13.54	330	110874	84.30	ng	0.00
44) 2-Fluorobiphenyl	11.21	172	755974	70.77	ng	0.00
78) Terphenyl-d14	17.01	244	446627	53.98	ng	0.00
Target Compounds						
49) Dimethylphthalate	11.90	163	163909	14.523	ng	Qvalue 99
70) Phenanthrene	14.68	178	120285	9.082	ng	97
74) Fluoranthene	16.45	202	86619	6.607	ng	98
77) Pyrene	16.76	202	93280	6.088	ng	97
80) Benzo(a)anthracene	18.34	228	36275	3.038	ng	95
82) Chrysene	18.38	228	38196m	3.201	ng	
87) Benzo(b)fluoranthene	19.62	252	27054m	2.805	ng	
89) Benzo(a)pyrene	19.96	252	22597	2.549	ng	# 91

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

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