

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060517\
 Data File : BF095693.D
 Acq On : 6 Jun 2017 1:54
 Operator : SJ/MA
 Sample : I3466-05
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-RO-6704-060217

Manual Integrations
 APPROVED

mohammad
 6/6/2017 5:13:07 PM

Quant Time: Jun 06 03:52:12 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	89645	20.00	ng	0.00
21) Naphthalene-d8	9.42	136	359695	20.00	ng	0.00
38) Acenaphthene-d10	12.25	164	153521	20.00	ng	0.00
63) Phenanthrene-d10	14.64	188	226041	20.00	ng	0.00
75) Chrysene-d12	18.35	240	181345	20.00	ng	0.00
86) Perylene-d12	20.02	264	138419	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	659980	117.31	ng	0.00
7) Phenol-d6	6.95	99	741842	110.15	ng	0.00
23) Nitrobenzene-d5	8.31	82	417379	72.06	ng	0.00
41) 2,4,6-Tribromophenol	13.54	330	121886	89.73	ng	0.00
44) 2-Fluorobiphenyl	11.21	172	705552	63.95	ng	0.00
78) Terphenyl-d14	17.01	244	417841	57.18	ng	0.00
Target Compounds						
49) Dimethylphthalate	11.90	163	170820	14.655	ng	Qvalue 99
70) Phenanthrene	14.68	178	83425	6.396	ng	94
71) Anthracene	14.77	178	27761	2.039	ng	97
74) Fluoranthene	16.45	202	104740	8.114	ng	98
77) Pyrene	16.76	202	114822	8.486	ng	97
80) Benzo(a)anthracene	18.34	228	49334	4.679	ng	96
82) Chrysene	18.39	228	46899	4.451	ng	96
87) Benzo(b)fluoranthene	19.62	252	42917m	4.952	ng	
89) Benzo(a)pyrene	19.97	252	30144	3.784	ng	# 93

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

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