

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122717\
 Data File : BF101784.D
 Acq On : 28 Dec 2017 1:02
 Operator : SJ/JU
 Sample : I6686-05 2X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-04-122217

Manual Integrations
 APPROVED

Sohil
 12/28/2017 2:03:19 PM

Quant Time: Dec 28 06:08:51 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 26 10:48:01 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	140503	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	530954	20.00	ng	0.00
38) Acenaphthene-d10	9.83	164	209222	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	332612	20.00	ng	0.00
75) Chrysene-d12	13.97	240	287657	20.00	ng	0.00
86) Perylene-d12	15.45	264	236812	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	486814	51.61	ng	0.01
7) Phenol-d6	6.44	99	614691	52.36	ng	0.00
23) Nitrobenzene-d5	7.36	82	349299	36.62	ng	0.00
41) 2,4,6-Tribromophenol	10.63	330	96914	48.07	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	619804	45.95	ng	0.00
78) Terphenyl-d14	12.90	244	439782	36.86	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.55	163	43089	2.560	ng	Qvalue # 98
51) Acenaphthene	9.86	154	30021	2.228	ng	99
70) Phenanthrene	11.35	178	93803	5.071	ng	99
74) Fluoranthene	12.55	202	154922	7.979	ng	97
77) Pyrene	12.79	202	137845	6.385	ng	100
80) Benzo(a)anthracene	13.97	228	68651	3.614	ng	98
82) Chrysene	14.00	228	67181	3.746	ng	98
87) Benzo(b)fluoranthene	15.03	252	66442m	4.603	ng	
89) Benzo(a)pyrene	15.39	252	43983	3.305	ng	# 86

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

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