

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120417\
 Data File : BF101091.D
 Acq On : 4 Dec 2017 12:15
 Operator : SJ/JU
 Sample : I6616-01RE
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P7-LOT-SP1-CRE

Manual Integrations
 APPROVED

Sohil
 12/5/2017 3:24:07 PM

Quant Time: Dec 04 15:46:12 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 04 13:07:49 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	53316	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	206981	20.00	ng	0.00
38) Acenaphthene-d10	9.88	164	98265	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	174968	20.00	ng	0.00
75) Chrysene-d12	14.02	240	122306	20.00	ng	0.00
86) Perylene-d12	15.49	264	126671	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	312273	96.78	ng	0.00
7) Phenol-d6	6.48	99	384341	98.11	ng	0.00
23) Nitrobenzene-d5	7.41	82	237578	68.47	ng	0.00
41) 2,4,6-Tribromophenol	10.67	330	103231	87.45	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	457768	63.74	ng	0.00
78) Terphenyl-d14	12.96	244	353881	63.29	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.59	163	26087	3.29	ng	99
74) Fluoranthene	12.59	202	29354	2.75	ng	94
77) Pyrene	12.82	202	26090	3.09	ng	99
82) Chrysene	14.04	228	18000m	2.51	ng	
83) Bis(2-ethylhexyl)phthalate	13.98	149	16261	3.07	ng	99
85) Indeno(1,2,3-cd)pyrene	16.96	276	13260	2.08	ng	90
87) Benzo(b)fluoranthene	15.06	252	27421m	3.61	ng	
89) Benzo(a)pyrene	15.42	252	17160	2.49	ng	# 89
91) Benzo(g,h,i)perylene	17.42	276	13123	2.29	ng	# 90

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

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