

Data Path : Z:\HPCHEM1\BNA F\DATA\BF041318\
 Data File : BF104476.D
 Acq On : 13 Apr 2018 14:14
 Operator : JU/SJ
 Sample : J2348-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 135-348

Manual Integrations
 APPROVED

Sohil
 4/16/2018 3:10:01 PM

Quant Time: Apr 13 17:52:37 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 13 17:38:31 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	142900	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	587245	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	276781	20.00	ng	0.00
63) Phenanthrene-d10	11.34	188	487776	20.00	ng	0.00
75) Chrysene-d12	13.99	240	323545	20.00	ng	0.00
86) Perylene-d12	15.45	264	284377	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	813590	87.06	ng	0.00
7) Phenol-d6	6.45	99	1003952	87.43	ng	0.00
23) Nitrobenzene-d5	7.37	82	631661	70.24	ng	0.00
41) 2,4,6-Tribromophenol	10.64	330	261612	91.12	ng	0.00
44) 2-Fluorobiphenyl	9.17	172	1029405	58.75	ng	0.00
78) Terphenyl-d14	12.93	244	977346	68.87	ng	0.00
Target Compounds						
10) Phenol	6.45	94	38198	3.135	ng	90
49) Dimethylphthalate	9.56	163	149025	6.828	ng	99
70) Phenanthrene	11.36	178	442090	16.275	ng	98
71) Anthracene	11.41	178	154160	5.583	ng	97
74) Fluoranthene	12.56	202	672094	23.681	ng	98
77) Pyrene	12.79	202	568138	23.690	ng	99
80) Benzo(a)anthracene	13.97	228	300551	15.549	ng	99
82) Chrysene	14.01	228	239369	12.057	ng	98
85) Indeno(1,2,3-cd)pyrene	16.90	276	115009m	6.819	ng	
87) Benzo(b)fluoranthene	15.03	252	263515m	14.672	ng	
88) Benzo(k)fluoranthene	15.05	252	83711m	4.937	ng	
89) Benzo(a)pyrene	15.39	252	195620	12.173	ng	96
90) Dibenzo(a,h)anthracene	16.92	278	31584	2.393	ng	96
91) Benzo(g,h,i)perylene	17.34	276	100372	7.849	ng	92

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

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