

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040618\
 Data File : BF104311.D
 Acq On : 6 Apr 2018 17:56
 Operator : JU/SJ
 Sample : J2239-01
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S-1-COMP

Manual Integrations
 APPROVED

Sohil
 4/9/2018 11:18:25 AM

Quant Time: Apr 06 18:39:17 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 06 16:44:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	149211	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	605478	20.00	ng	0.00
38) Acenaphthene-d10	9.84	164	276751	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	450321	20.00	ng	0.00
75) Chrysene-d12	13.96	240	318773	20.00	ng	-0.01
86) Perylene-d12	15.42	264	327992	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	1295571	132.77	ng	0.01
7) Phenol-d6	6.43	99	1602059	133.62	ng	0.00
23) Nitrobenzene-d5	7.37	82	960422	103.58	ng	0.00
41) 2,4,6-Tribromophenol	10.63	330	357362	124.48	ng	0.00
44) 2-Fluorobiphenyl	9.16	172	1498637	85.55	ng	0.00
78) Terphenyl-d14	12.92	244	1144252	81.84	ng	0.00
Target Compounds						
10) Phenol	6.45	94	122773	9.651	ng	97
49) Dimethylphthalate	9.56	163	283696	12.999	ng	99
56) 2,4-Dinitrotoluene	10.02	165	16070	3.034	ng	# 96
70) Phenanthrene	11.35	178	63364	2.527	ng	99
74) Fluoranthene	12.53	202	143093	5.461	ng	98
77) Pyrene	12.76	202	126947	5.373	ng	99
80) Benzo(a)anthracene	13.95	228	65310	3.429	ng	94
82) Chrysene	13.99	228	67646	3.458	ng	98
85) Indeno(1,2,3-cd)pyrene	16.84	276	46794	2.816	ng	94
87) Benzo(b)fluoranthene	15.00	252	106876m	5.159	ng	
89) Benzo(a)pyrene	15.36	252	70066	3.780	ng	97
91) Benzo(a,h,i)perylene	17.27	276	43844	2.973	ng	98

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

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