

Data Path : Z:\HPCHEM1\BNA F\DATA\BF041618\
 Data File : BF104549.D
 Acq On : 16 Apr 2018 22:40
 Operator : JU/SJ
 Sample : J2370-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CG-RO-721194

Manual Integrations
 APPROVED

Sohil
 4/17/2018 10:09:19 AM

Quant Time: Apr 17 02:01:10 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 16 12:15:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	153586	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	605067	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	244995	20.00	ng	0.00
63) Phenanthrene-d10	11.34	188	363169	20.00	ng	0.00
75) Chrysene-d12	13.99	240	288563	20.00	ng	0.00
86) Perylene-d12	15.46	264	226970	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	710075	70.69	ng	0.01
7) Phenol-d6	6.45	99	903340	73.20	ng	0.00
23) Nitrobenzene-d5	7.37	82	564140	60.88	ng	0.00
41) 2,4,6-Tribromophenol	10.65	330	157984	62.16	ng	0.00
44) 2-Fluorobiphenyl	9.17	172	1004126	64.75	ng	0.00
78) Terphenyl-d14	12.93	244	705220	55.72	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.56	163	105879	5.480	ng	100
70) Phenanthrene	11.36	178	204871	10.130	ng	99
74) Fluoranthene	12.56	202	142616	6.749	ng	99
77) Pyrene	12.79	202	239514	11.198	ng	99
80) Benzo(a)anthracene	13.98	228	81515m	4.729	ng	
82) Chrysene	14.02	228	79253	4.476	ng	97
87) Benzo(b)fluoranthene	15.03	252	52930m	3.692	ng	
89) Benzo(a)pyrene	15.39	252	46120	3.596	ng	# 93
91) Benzo(g,h,i)perylene	17.35	276	20792	2.037	ng	95

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

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