

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012017\
 Data File : BF092372.D
 Acq On : 20 Jan 2017 14:52
 Operator : UM/SJ
 Sample : I1304-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ROLL-OFF-15-41-COMP

Manual Integrations
 APPROVED

mohammad
 1/23/2017 2:44:44 PM

Quant Time: Jan 20 23:04:37 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 20 22:22:14 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.52	152	139809	20.00	ng	0.00
21) Naphthalene-d8	7.81	136	849639	20.00	ng	0.00
38) Acenaphthene-d10	9.56	164	444263	20.00	ng	0.00
63) Phenanthrene-d10	11.03	188	670858	20.00	ng	0.00
75) Chrysene-d12	13.66	240	438321	20.00	ng	0.00
86) Perylene-d12	15.01	264	364678	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.11	112	1368146	163.40	ng	0.02
7) Phenol-d6	6.19	99	1854662	181.42	ng	0.00
23) Nitrobenzene-d5	7.10	82	1292865	84.61	ng	0.00
41) 2,4,6-Tribromophenol	10.35	330	464635	126.38	ng	0.00
44) 2-Fluorobiphenyl	8.90	172	2292738	109.45	ng	0.00
78) Terphenyl-d14	12.62	244	1670837	92.45	ng	0.00
Target Compounds						
10) Phenol	6.20	94	56330	4.84	ng	Qvalue # 26
49) Dimethylphthalate	9.30	163	547028	19.25	ng	100
77) Pyrene	12.46	202	145890	4.54	ng	97
80) Benzo(a)anthracene	13.65	228	77502m	3.13	ng	
82) Chrysene	13.69	228	86020m	3.47	ng	
85) Indeno(1,2,3-cd)pyrene	16.23	276	56860	2.64	ng	93
87) Benzo(b)fluoranthene	14.66	252	183219m	8.07	ng	
88) Benzo(k)fluoranthene	14.67	252	57966m	2.99	ng	
89) Benzo(a)pyrene	14.97	252	101145	5.02	ng	97
91) Benzo(g,h,i)perylene	16.59	276	50577	2.94	ng	# 94

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

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