

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010319\
 Data File : BF111838.D
 Acq On : 3 Jan 2019 23:05
 Operator : JU/SJ
 Sample : K1019-03
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEST-PIT-D-11-12-B

Manual Integrations
 APPROVED

Sohil
 1/4/2019 9:56:54 AM

Quant Time: Jan 04 04:02:30 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 19 17:39:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	240893	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	889266	20.00	ng	0.00
39) Acenaphthene-d10	9.94	164	412337	20.00	ng	0.00
64) Phenanthrene-d10	11.43	188	601419	20.00	ng	0.01
76) Chrysene-d12	14.07	240	458319	20.00	ng	0.02
87) Perylene-d12	15.56	264	315509	20.00	ng	0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	1338481	94.71	ng	0.02
7) Phenol-d6	6.55	99	1790013	99.52	ng	0.02
23) Nitrobenzene-d5	7.46	82	1089360	71.31	ng	0.00
42) 2,4,6-Tribromophenol	10.73	330	347907	71.41	ng	0.01
45) 2-Fluorobiphenyl	9.26	172	2018113	71.34	ng	0.00
79) Terphenyl-d14	13.01	244	1483031	61.36	ng	0.01
Target Compounds						
10) Phenol	6.56	94	82102	4.046	ng	96
50) Dimethylphthalate	9.65	163	329553	10.929	ng	98
58) Fluorene	10.49	166	62988	2.330	ng	100
71) Phenanthrene	11.45	178	614823	19.276	ng	94
72) Anthracene	11.50	178	73225	2.287	ng	93
75) Fluoranthene	12.64	202	551474	17.074	ng	96
78) Pyrene	12.87	202	434573	13.427	ng	95
81) Benzo(a)anthracene	14.06	228	172807	6.607	ng	97
83) Chrysene	14.09	228	158280	6.157	ng	99
86) Indeno(1,2,3-cd)pyrene	17.05	276	46944	2.148	ng	# 87
88) Benzo(b)fluoranthene	15.12	252	137053m	7.433	ng	
89) Benzo(k)fluoranthene	15.14	252	49708m	2.832	ng	
90) Benzo(a)pyrene	15.49	252	89536	5.345	ng	97
92) Benzo(a,h,i)perylene	17.50	276	42845	2.991	ng	# 91

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

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