

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030419\
 Data File : BF112818.D
 Acq On : 4 Mar 2019 12:29
 Operator : JU/SJ
 Sample : K1670-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-02-032-A

Manual Integrations
 APPROVED

Sohil
 3/5/2019 11:00:34 AM

Quant Time: Mar 05 01:01:36 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 01 15:40:43 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	185039	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	715858	20.00	ng	0.00
39) Acenaphthene-d10	9.77	164	328228	20.00	ng	0.00
64) Phenanthrene-d10	11.25	188	607034	20.00	ng	0.00
76) Chrysene-d12	13.88	240	443868	20.00	ng	0.00
87) Perylene-d12	15.29	264	471811	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	1032966	86.84	ng	0.01
7) Phenol-d6	6.38	99	1340408	89.70	ng	0.01
23) Nitrobenzene-d5	7.30	82	820626	68.60	ng	0.00
42) 2,4,6-Tribromophenol	10.56	330	371602	106.64	ng	0.00
45) 2-Fluorobiphenyl	9.10	172	1450491	69.86	ng	0.00
79) Terphenyl-d14	12.83	244	1254029	54.77	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.49	163	127926	5.286	ng	Qvalue 100
71) Phenanthrene	11.27	178	126555	4.190	ng	98
75) Fluoranthene	12.46	202	370113	11.438	ng	97
78) Pyrene	12.69	202	319684	8.543	ng	98
81) Benzo(a)anthracene	13.87	228	215841	7.016	ng	96
83) Chrysene	13.90	228	208055	6.904	ng	96
86) Indeno(1,2,3-cd)pyrene	16.65	276	90360	3.517	ng	# 94
88) Benzo(b)fluoranthene	14.89	252	254945m	8.354	ng	
89) Benzo(k)fluoranthene	14.92	252	92415m	3.343	ng	
90) Benzo(a)pyrene	15.23	252	167029	6.188	ng	99
92) Benzo(g,h,i)perylene	17.06	276	71635	3.097	ng	95

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

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