

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF030119\
 Data File : BF112804.D
 Acq On : 1 Mar 2019 17:16
 Operator : JU/SJ
 Sample : K1670-03 2X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 3/4/2019 4:55:31 PM

Quant Time: Mar 01 23:21:09 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	169591	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	623310	20.00	ng	0.00
39) Acenaphthene-d10	9.77	164	248073	20.00	ng	0.00
64) Phenanthrene-d10	11.25	188	420330	20.00	ng	0.00
76) Chrysene-d12	13.87	240	434795	20.00	ng	0.00
87) Perylene-d12	15.29	264	378676	20.00	ng	0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	494676	45.37	ng	0.00
7) Phenol-d6	6.37	99	627791	45.84	ng	0.00
23) Nitrobenzene-d5	7.30	82	354772	34.06	ng	0.00
42) 2,4,6-Tribromophenol	10.56	330	126360	47.98	ng	0.00
45) 2-Fluorobiphenyl	9.10	172	600194	34.00	ng	0.00
79) Terphenyl-d14	12.83	244	536772	23.93	ng	0.00

Target Compounds				Qvalue		
10) Phenol	6.38	94	44826	2.805	ng	92
49) Acenaphthylene	9.63	152	61099	2.278	ng	98
50) Dimethylphthalate	9.49	163	43439	2.375	ng	# 96
75) Fluoranthene	12.46	202	210013	9.373	ng	94
78) Pyrene	12.68	202	244822	6.679	ng	98
81) Benzo(a)anthracene	13.86	228	213499m	7.085	ng	
83) Chrysene	13.90	228	184998	6.267	ng	96
86) Indeno(1,2,3-cd)pyrene	16.64	276	91996m	3.656	ng	
88) Benzo(b)fluoranthene	14.89	252	308842m	12.609	ng	
89) Benzo(k)fluoranthene	14.91	252	69942m	3.152	ng	
90) Benzo(a)pyrene	15.23	252	179644	8.292	ng	97
92) Benzo(a,h,i)perylene	17.04	276	82758	4.458	ng	97

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

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