

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031319\
 Data File : BF113042.D
 Acq On : 13 Mar 2019 21:44
 Operator : JU/SJ
 Sample : K1870-01 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-1

Manual Integrations
 APPROVED

Sohil
 3/14/2019 10:50:10 AM

Quant Time: Mar 14 02:36:13 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 13 03:42:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	153163	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	573330	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	318052	20.00	ng	-0.01
64) Phenanthrene-d10	11.23	188	655092	20.00	ng	-0.01
76) Chrysene-d12	13.86	240	523564	20.00	ng	-0.01
87) Perylene-d12	15.26	264	505946	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	424740	43.14	ng	-0.01
7) Phenol-d6	6.37	99	582230	47.07	ng	0.00
23) Nitrobenzene-d5	7.28	82	332968	34.75	ng	-0.02
42) 2,4,6-Tribromophenol	10.54	330	206633	61.20	ng	-0.01
45) 2-Fluorobiphenyl	9.08	172	654424	28.25	ng	-0.01
79) Terphenyl-d14	12.81	244	667121	24.70	ng	-0.01
Target Compounds						Qvalue
50) Dimethylphthalate	9.47	163	62044	2.646	ng	98
71) Phenanthrene	11.26	178	72361	2.220	ng	97
75) Fluoranthene	12.44	202	162285	4.647	ng	94
78) Pyrene	12.66	202	154634	3.503	ng	98
81) Benzo(a)anthracene	13.84	228	76784	2.116	ng	99
83) Chrysene	13.88	228	75403	2.121	ng	96
88) Benzo(b)fluoranthene	14.87	252	108901m	3.328	ng	
90) Benzo(a)pyrene	15.20	252	77546	2.679	ng	97

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

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