

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP101819\
 Data File : BP000868.D
 Acq On : 18 Oct 2019 05:05
 Operator : JU
 Sample : K5394-01
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 T-17-F1-(0-30)

Manual Integrations
 APPROVED

mohammad
 10/20/2019 8:16:55 PM

Quant Time: Oct 18 08:43:25 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP100119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 01 18:03:42 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	200908	20.00	ng	-0.02
21) Naphthalene-d8	8.03	136	671863	20.00	ng	-0.02
39) Acenaphthene-d10	9.80	164	347343	20.00	ng	-0.02
64) Phenanthrene-d10	11.30	188	601736	20.00	ng	-0.01
76) Chrysene-d12	13.97	240	504208	20.00	ng	0.00
87) Perylene-d12	15.46	264	561134	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	1060440	84.84	ng	-0.01
7) Phenol-d6	6.40	99	1319225	87.59	ng	-0.01
23) Nitrobenzene-d5	7.31	82	723397	65.76	ng	-0.02
42) 2,4,6-Tribromophenol	10.59	330	367674	99.34	ng	-0.02
45) 2-Fluorobiphenyl	9.12	172	1520067	70.81	ng	-0.02
79) Terphenyl-d14	12.90	244	1567286	62.94	ng	-0.01
Target Compounds						Qvalue
10) Phenol	6.41	94	58434	3.789	ng	98
31) Naphthalene	8.05	128	251782	8.024	ng	99
37) 2-Methylnaphthalene	8.75	142	109633	5.205	ng	98
38) 1-Methylnaphthalene	8.85	142	84182m	4.230	ng	
49) Acenaphthylene	9.66	152	176613	5.774	ng	99
50) Dimethylphthalate	9.51	163	305218	12.646	ng	99
71) Phenanthrene	11.32	178	629371	20.826	ng	100
72) Anthracene	11.37	178	169413	5.654	ng	99
75) Fluoranthene	12.53	202	821164	24.194	ng	99
78) Pyrene	12.76	202	1004274	28.879	ng	99
81) Benzo(a)anthracene	13.97	228	540991	17.585	ng	# 90
83) Chrysene	14.00	228	557703	18.470	ng	98
86) Indeno(1,2,3-cd)pyrene	16.93	276	309734	8.212	ng	# 94
88) Benzo(b)fluoranthene	15.03	252	755200m	23.723	ng	
89) Benzo(k)fluoranthene	15.06	252	173468m	5.724	ng	
90) Benzo(a)pyrene	15.40	252	527434	17.773	ng	96
91) Dibenzo(a,h)anthracene	16.94	278	84969m	2.952	ng	
92) Benzo(g,h,i)perylene	17.38	276	320950	10.974	ng	98

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

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