

Data Path : Z:\HPCHEM1\BNA F\DATA\BF071917\
 Data File : BF096883.D
 Acq On : 19 Jul 2017 21:24
 Operator : SJ/JU
 Sample : I4280-04
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-01-071817-B

Manual Integrations
 APPROVED

Sohil
 7/20/2017 3:05:03 PM

Quant Time: Jul 20 05:07:02 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF071317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 18 13:33:22 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.59	152	116281	20.00	ng	-0.01
21) Naphthalene-d8	7.87	136	503005	20.00	ng	-0.02
38) Acenaphthene-d10	9.62	164	178557	20.00	ng	-0.02
63) Phenanthrene-d10	11.10	188	254757	20.00	ng	0.00
75) Chrysene-d12	13.73	240	199689	20.00	ng	0.00
86) Perylene-d12	15.10	264	165586	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.20	112	680125	87.94	ng	0.00
7) Phenol-d6	6.24	99	832234	89.68	ng	-0.01
23) Nitrobenzene-d5	7.15	82	485169	59.74	ng	-0.02
41) 2,4,6-Tribromophenol	10.40	330	127557	83.69	ng	-0.02
44) 2-Fluorobiphenyl	8.95	172	804721	71.20	ng	-0.01
78) Terphenyl-d14	12.68	244	441988	38.39	ng	-0.01
Target Compounds						Qvalue
10) Phenol	6.25	94	33409	3.184	ng	94
20) 3+4-Methylphenols	7.01	107	46066	5.851	ng	# 57
31) Naphthalene	7.89	128	284286	11.931	ng	99
37) 2-Methylnaphthalene	8.58	142	197050	12.971	ng	98
48) Acenaphthylene	9.47	152	121258	6.749	ng	# 96
49) Dimethylphthalate	9.35	163	96364	7.167	ng	100
51) Acenaphthene	9.65	154	108541	10.226	ng	96
54) Dibenzofuran	9.82	168	98431	6.618	ng	# 95
57) Fluorene	10.16	166	266947	24.288	ng	100
70) Phenanthrene	11.13	178	1050482	75.836	ng	99
71) Anthracene	11.17	178	329173	23.503	ng	99
72) Carbazole	11.33	167	90857	6.699	ng	98
74) Fluoranthene	12.31	202	832099	62.758	ng	99
77) Pyrene	12.53	202	925516	51.069	ng	99
80) Benzo(a)anthracene	13.72	228	399559	32.113	ng	95
82) Chrysene	13.75	228	341014m	27.841	ng	
85) Indeno(1,2,3-cd)pyrene	16.37	276	109048	12.473	ng	93
87) Benzo(b)fluoranthene	14.73	252	317932m	31.835	ng	
88) Benzo(k)fluoranthene	14.74	252	100800m	10.659	ng	
89) Benzo(a)pyrene	15.04	252	257112	28.070	ng	99
90) Dibenzo(a,h)anthracene	16.38	278	33344	3.956	ng	# 71
91) Benzo(g,h,i)perylene	16.76	276	111269	12.329	ng	98

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

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