

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121416\
 Data File : BF091594.D
 Acq On : 14 Dec 2016 19:51
 Operator : UM/SJ
 Sample : H6006-08
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-8-0-5

Manual Integrations
 APPROVED

sohil
 12/15/2016 5:40:11 PM

Quant Time: Dec 15 01:19:00 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 09 19:00:21 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	308380	20.00	ng	-0.02
21) Naphthalene-d8	8.08	136	1239182	20.00	ng	-0.02
38) Acenaphthene-d10	9.84	164	657191	20.00	ng	-0.02
63) Phenanthrene-d10	11.31	188	1006465	20.00	ng	-0.02
75) Chrysene-d12	13.95	240	689345	20.00	ng	-0.01
86) Perylene-d12	15.36	264	634504	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	934442	51.05	ng	-0.02
7) Phenol-d6	6.43	99	1382726	60.59	ng	-0.02
23) Nitrobenzene-d5	7.36	82	969902	43.68	ng	-0.02
41) 2,4,6-Tribromophenol	10.62	330	573661	105.09	ng	-0.01
44) 2-Fluorobiphenyl	9.16	172	2028874	53.34	ng	-0.02
78) Terphenyl-d14	12.90	244	1910955	62.90	ng	-0.02
Target Compounds						
10) Phenol	6.44	94	153552	5.76	ng	82
49) Dimethylphthalate	9.55	163	598440	14.38	ng	98
51) Acenaphthene	9.86	154	100126	2.57	ng	96
54) Dibenzofuran	10.03	168	133762	2.77	ng	96
70) Phenanthrene	11.34	178	2085938	41.85	ng	99
71) Anthracene	11.39	178	662256	12.32	ng	98
72) Carbazole	11.54	167	176750	3.75	ng	99
74) Fluoranthene	12.52	202	1967049	37.69	ng	94
77) Pyrene	12.75	202	1682140	30.76	ng	98
80) Benzo(a)anthracene	13.94	228	841033	21.01	ng	99
82) Chrysene	13.97	228	698946m	16.69	ng	
83) Bis(2-ethylhexyl)phthalate	13.94	149	160027	5.85	ng	97
85) Indeno(1,2,3-cd)pyrene	16.72	276	336923	8.94	ng	95
87) Benzo(b)fluoranthene	14.96	252	705389m	18.65	ng	
88) Benzo(k)fluoranthene	14.97	252	330532m	9.94	ng	
89) Benzo(a)pyrene	15.30	252	552276	16.21	ng	99
90) Dibenzo(a,h)anthracene	16.73	278	87191	2.86	ng	94
91) Benzo(g,h,i)perylene	17.13	276	318675	10.26	ng	97

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

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