

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF110216\
 Data File : BF091000.D
 Acq On : 3 Nov 2016 00:34
 Operator : UM/SJ
 Sample : H5493-25DL 10X
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleID :
 SB-6(0-2)DL

Manual Integrations
 APPROVED

Umangi
 11/3/2016 7:14:10 PM

Quant Time: Nov 03 02:15:35 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF102616.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 01 11:53:50 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	161182	20.00	ng	-0.02
21) Naphthalene-d8	8.01	136	647026	20.00	ng	-0.03
38) Acenaphthene-d10	9.77	164	332284	20.00	ng	-0.02
63) Phenanthrene-d10	11.25	188	458043	20.00	ng	-0.01
75) Chrysene-d12	13.92	240	351079	20.00	ng	0.01
86) Perylene-d12	15.32	264	396537	20.00	ng	0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.32	112	1263767	137.09	ng	-0.01
7) Phenol-d6	6.37	99	1511556	121.53	ng	-0.02
23) Nitrobenzene-d5	7.29	82	968504	97.99	ng	-0.03
41) 2,4,6-Tribromophenol	10.56	330	299056	87.58	ng	-0.02
44) 2-Fluorobiphenyl	9.09	172	1510350	79.82	ng	-0.02
78) Terphenyl-d14	12.84	244	1148201	82.40	ng	-0.01

Target Compounds

						Qvalue
31) Naphthalene	8.03	128	612420	20.25	ng	98
37) 2-Methylnaphthalene	8.73	142	153107	7.84	ng	# 90
45) 1,1'-Biphenyl	9.20	154	70236	2.93	ng	94
48) Acenaphthylene	9.63	152	1845898	66.96	ng	# 93
49) Dimethylphthalate	9.49	163	568275	25.60	ng	# 98
51) Acenaphthene	9.80	154	499164	26.29	ng	87
54) Dibenzofuran	9.97	168	461147	18.82	ng	# 76
57) Fluorene	10.32	166	1296613	64.61	ng	91
70) Phenanthrene	11.29	178	4846243	207.76	ng	# 91
71) Anthracene	11.33	178	2668796	108.66	ng	93
72) Carbazole	11.49	167	219657	10.13	ng	# 85
73) Di-n-butylphthalate	11.83	149	87469	3.13	ng	# 80
74) Fluoranthene	12.50	202	6491079	254.11	ng	89
77) Pyrene	12.72	202	5665289m	254.99	ng	
80) Benzo(a)anthracene	13.91	228	4122706	221.27	ng	# 82
82) Chrysene	13.94	228	3618846m	192.02	ng	
85) Indeno(1,2,3-cd)pyrene	16.68	276	2230458m	133.55	ng	
87) Benzo(b)fluoranthene	14.93	252	4774056m	179.04	ng	
88) Benzo(k)fluoranthene	14.96	252	1396493m	66.27	ng	
89) Benzo(a)pyrene	15.28	252	3479822	151.89	ng	# 83
90) Dibenzo(a,h)anthracene	16.68	278	625109	32.23	ng	# 77
91) Benzo(g,h,i)perylene	17.09	276	2013357	110.22	ng	# 73

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

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