

Data Path : Z:\HPCHEM1\BNA F\DATA\BF110116\
 Data File : BF090972.D
 Acq On : 2 Nov 2016 1:00
 Operator : UM/SJ
 Sample : H5502-23
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-90-0.0-1.0

Manual Integrations
 APPROVED

Umangi
 11/2/2016 5:42:44 PM

Quant Time: Nov 02 02:31:09 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	6.74	152	177993	20.00	ng	-0.01
21) Naphthalene-d8	8.03	136	690649	20.00	ng	-0.01
38) Acenaphthene-d10	9.78	164	333187	20.00	ng	-0.01
63) Phenanthrene-d10	11.28	188	463547	20.00	ng	0.01
75) Chrysene-d12	13.93	240	272097	20.00	ng	0.02
86) Perylene-d12	15.33	264	382350	20.00	ng	0.03

System Monitoring Compounds

5) 2-Fluorophenol	5.33	112	1055208	103.65	ng	0.00
7) Phenol-d6	6.38	99	1577044	114.82	ng	-0.01
23) Nitrobenzene-d5	7.31	82	1056646	100.16	ng	-0.01
41) 2,4,6-Tribromophenol	10.58	330	262926	76.79	ng	0.00
44) 2-Fluorobiphenyl	9.10	172	1379963	72.73	ng	-0.01
78) Terphenyl-d14	12.85	244	1163305	107.71	ng	0.00

Target Compounds

						Qvalue
6) Aniline	6.43	93	135081	8.61	ng	# 58
10) Phenol	6.40	94	87500	5.78	ng	# 1
22) Acetophenone	7.15	105	47224	3.32	ng	# 82
31) Naphthalene	8.05	128	478050	14.81	ng	# 99
37) 2-Methylnaphthalene	8.74	142	169843	8.15	ng	# 89
45) 1,1'-Biphenyl	9.21	154	98224	4.09	ng	# 96
48) Acenaphthylene	9.64	152	261337	9.45	ng	# 96
49) Dimethylphthalate	9.50	163	463001	20.80	ng	# 97
51) Acenaphthene	9.81	154	649316	34.11	ng	# 88
54) Dibenzofuran	9.98	168	791580	32.22	ng	# 84
57) Fluorene	10.33	166	822862	40.89	ng	# 92
70) Phenanthrene	11.31	178	4564691m	193.37	ng	# 95
71) Anthracene	11.36	178	1851754	74.50	ng	# 95
72) Carbazole	11.50	167	874775	39.85	ng	# 92
74) Fluoranthene	12.50	202	6377044	246.69	ng	# 90
77) Pyrene	12.74	202	5268267	305.95	ng	# 90
80) Benzo(a)anthracene	13.92	228	4469308	309.50	ng	# 87
82) Chrysene	13.96	228	2598743m	177.92	ng	# 96
83) Bis(2-ethylhexyl)phthalate	13.91	149	107089	10.53	ng	# 96
85) Indeno(1,2,3-cd)pyrene	16.71	276	1931638	149.23	ng	# 86
87) Benzo(b)fluoranthene	14.95	252	4752128m	184.84	ng	# 84
88) Benzo(k)fluoranthene	14.97	252	1075481m	52.93	ng	# 80
89) Benzo(a)pyrene	15.29	252	2996214	135.63	ng	# 84
90) Dibenz(a,h)anthracene	16.71	278	592285	31.67	ng	# 80
91) Benzo(g,h,i)perylene	17.12	276	1590925	90.32	ng	# 77

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

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