

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121516\
 Data File : BF091671.D
 Acq On : 16 Dec 2016 12:31
 Operator : UM/SJ
 Sample : H6007-12DL 2X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-6-0-5DL

Quant Time: Dec 16 17:06:07 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 16 00:35:41 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.78	152	173642	20.00	ng	0.00
21) Naphthalene-d8	8.06	136	616425	20.00	ng	-0.01
38) Acenaphthene-d10	9.82	164	289988	20.00	ng	0.00
63) Phenanthrene-d10	11.30	188	384405	20.00	ng	-0.01
75) Chrysene-d12	13.94	240	330752	20.00	ng	0.00
86) Perylene-d12	15.35	264	257230	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	592112	57.45	ng	0.00
7) Phenol-d6	6.42	99	701940	54.63	ng	-0.01
23) Nitrobenzene-d5	7.34	82	466139	42.21	ng	-0.01
41) 2,4,6-Tribromophenol	10.61	330	134040	55.65	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	769306	45.84	ng	0.00
78) Terphenyl-d14	12.89	244	510688	35.04	ng	-0.01

Target Compounds						Qvalue
19) n-Nitroso-di-n-propylamine	7.34	70	64514	8.03	ng	# 68
20) 3+4-Methylphenols	7.21	107	1314	Below	Cal	# 81
25) Isophorone	7.34	82	410488	20.91	ng	# 67
32) Benzoic acid	7.87	122	648	4.99	ng	# 1
48) Acenaphthylene	9.69	152	68865	2.78	ng	97
49) Dimethylphthalate	9.54	163	163204	8.89	ng	# 97
51) Acenaphthene	9.86	154	65263	3.79	ng	97
55) 4-Nitrophenol	10.03	139	15533	4.41	ng	# 2
56) 2,4-Dinitrotoluene	9.82	165	39831	7.88	ng	# 34
57) Fluorene	10.37	166	48382	Below	Cal	99
64) 4,6-Dinitro-2-methylphenol	10.61	198	438	3.12	ng	# 1
66) 4-Bromophenyl-phenylether	10.61	248	8758	2.22	ng	# 1
70) Phenanthrene	11.33	178	729164	38.30	ng	99
71) Anthracene	11.38	178	197428	9.62	ng	97
72) Carbazole	11.54	167	75992	4.22	ng	97
74) Fluoranthene	12.52	202	1138140	57.10	ng	97
77) Pyrene	12.74	202	1064251	40.56	ng	97
80) Benzo(a)anthracene	13.93	228	683346	35.58	ng	97
82) Chrysene	13.96	228	494600	24.61	ng	97
85) Indeno(1,2,3-cd)pyrene	16.71	276	227005	12.55	ng	93
87) Benzo(b)fluoranthene	14.96	252	848445	55.33	ng	100
88) Benzo(k)fluoranthene	14.96	252	848445	62.93	ng	98
89) Benzo(a)pyrene	15.29	252	426328	30.86	ng	99
90) Dibenzo(a,h)anthracene	16.88	278	52626	4.25	ng	96
91) Benzo(g,h,i)perylene	17.12	276	213518	16.95	ng	99

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

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