

Data Path : Z:\HPCHEM1\BNA G\DATA\BG030216\
 Data File : BG021212.D
 Acq On : 2 Mar 2016 14:25
 Operator : UM/SJ
 Sample : H1578-03DL 25X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CEB-7(2.5-5)20160224DL

Manual Integrations
 APPROVED

Sohil
 3/3/2016 8:22:52 AM

Quant Time: Mar 02 15:10:27 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 02 14:51:49 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.14	152	11888	20.00	ng	0.00
21) Naphthalene-d8	10.95	136	66606	20.00	ng	0.00
38) Acenaphthene-d10	14.76	164	46165	20.00	ng	0.00
63) Phenanthrene-d10	17.49	188	108010	20.00	ng	0.00
75) Chrysene-d12	21.77	240	110148	20.00	ng	0.00
86) Perylene-d12	25.05	264	103306	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.70	112	3175	4.25	ng	0.00
7) Phenol-d6	7.30	99	5058	4.11	ng	0.00
23) Nitrobenzene-d5	9.31	82	3743	2.38	ng	0.00
41) 2,4,6-Tribromophenol	16.24	330	1389	2.89	ng	0.00
44) 2-Fluorobiphenyl	13.38	172	6671	2.05	ng	0.00
78) Terphenyl-d14	20.09	244	9063	1.99	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	11.01	128	8123	2.34	ng	99
51) Acenaphthene	14.82	154	6245	2.04	ng	95
57) Fluorene	15.80	166	9422	2.42	ng	93
70) Phenanthrene	17.54	178	136503	23.42	ng	98
71) Anthracene	17.63	178	20761	3.48	ng	99
74) Fluoranthene	19.54	202	136824	19.98	ng	97
77) Pyrene	19.90	202	138812	21.12	ng	99
80) Benzo(a)anthracene	21.75	228	59039	9.05	ng	97
82) Chrysene	21.81	228	62227	10.07	ng	99
85) Indeno(1,2,3-cd)pyrene	28.78	276	29342	4.14	ng	# 92
87) Benzo(b)fluoranthene	24.00	252	61281m	9.43	ng	
88) Benzo(k)fluoranthene	24.06	252	22195m	3.41	ng	
89) Benzo(a)pyrene	24.89	252	50115	7.98	ng	93
91) Benzo(a,h,i)perylene	29.95	276	28077	4.68	ng	96

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

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