

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040416\
 Data File : BF086053.D
 Acq On : 4 Apr 2016 22:38
 Operator : UM/SJ
 Sample : H2165-02 10X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEC-B2COMP

Manual Integrations
 APPROVED

Sohil
 4/5/2016 6:39:10 PM

Quant Time: Apr 05 03:41:38 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 01 23:17:06 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.76	152	10947	20.00	ng	-0.01
21) Naphthalene-d8	8.05	136	40126	20.00	ng	-0.01
38) Acenaphthene-d10	9.80	164	17499	20.00	ng	-0.01
63) Phenanthrene-d10	11.29	188	26675	20.00	ng	-0.01
75) Chrysene-d12	13.92	240	20924	20.00	ng	-0.01
86) Perylene-d12	15.32	264	16888	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	75885	118.49	ng	0.00
7) Phenol-d6	6.42	99	100848	123.97	ng	0.00
23) Nitrobenzene-d5	7.33	82	54809	80.55	ng	-0.01
41) 2,4,6-Tribromophenol	10.59	330	15150	92.24	ng	-0.01
44) 2-Fluorobiphenyl	9.13	172	100990	95.96	ng	-0.01
78) Terphenyl-d14	12.87	244	69368	77.93	ng	-0.01
Target Compounds						Qvalue
31) Naphthalene	8.08	128	28172	13.96	ng	97
37) 2-Methylnaphthalene	8.76	142	7782	5.90	ng	96
48) Acenaphthylene	9.66	152	4302	2.46	ng	99
49) Dimethylphthalate	9.53	163	10742	8.28	ng	97
51) Acenaphthene	9.84	154	9024	8.66	ng	96
54) Dibenzofuran	10.01	168	10788	7.84	ng	97
57) Fluorene	10.35	166	13957	11.87	ng	96
70) Phenanthrene	11.31	178	85420	58.70	ng	97
71) Anthracene	11.36	178	26803	17.58	ng	98
72) Carbazole	11.53	167	10166	7.76	ng	97
74) Fluoranthene	12.50	202	85039	56.18	ng	90
77) Pvrene	12.73	202	70350	47.71	ng	99
79) Butylbenzylphthalate	13.35	149	3243	4.88	ng	# 70
80) Benzo(a)anthracene	13.91	228	33737	27.35	ng	94
82) Chrysene	13.94	228	24588	20.70	ng	# 93
85) Indeno(1,2,3-cd)pyrene	16.68	276	11102	11.52	ng	98
87) Benzo(b)fluoranthene	14.93	252	33004m	31.07	ng	
88) Benzo(k)fluoranthene	14.96	252	13852m	14.72	ng	
91) Benzo(a,h,i)perylene	17.11	276	11138	15.20	ng	# 79

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

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