

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040816\
 Data File : BF086207.D
 Acq On : 9 Apr 2016 9:30
 Operator : UM/SJ
 Sample : H2310-06DL 150X
 Misc :
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PEDBR-SB4-0-6.0DL

Manual Integrations
 APPROVED

Sohil
 4/11/2016 5:45:34 PM

Quant Time: Apr 11 17:18:57 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	6.73	152	101457	20.00	ng	-0.05
21) Naphthalene-d8	8.01	136	390356	20.00	ng	-0.06
38) Acenaphthene-d10	9.77	164	158514	20.00	ng	-0.05
63) Phenanthrene-d10	11.24	188	259715	20.00	ng	-0.06
75) Chrysene-d12	13.88	240	224368	20.00	ng	-0.05
86) Perylene-d12	15.28	264	178075	20.00	ng	-0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	1879	0.32	ng	0.03
7) Phenol-d6	6.43	99	5052	0.67	ng	0.01
23) Nitrobenzene-d5	7.33	82	2483	0.38	ng	-0.01
41) 2,4,6-Tribromophenol	10.57	330	374	0.25	ng	-0.03
44) 2-Fluorobiphenyl	9.09	172	5728	0.60	ng	-0.05
78) Terphenyl-d14	12.83	244	3474	0.36	ng	-0.05
Target Compounds						Qvalue
31) Naphthalene	8.03	128	50613	2.58	ng	99
51) Acenaphthene	9.79	154	43866	4.65	ng	99
54) Dibenzofuran	9.97	168	42675	3.42	ng	98
57) Fluorene	10.30	166	38087	3.58	ng	98
70) Phenanthrene	11.26	178	425602	30.04	ng	99
71) Anthracene	11.32	178	133031	8.96	ng	97
72) Carbazole	11.48	167	30690	2.41	ng	98
74) Fluoranthene	12.45	202	368410	25.00	ng	96
77) Pyrene	12.68	202	330676	20.91	ng	100
80) Benzo(a)anthracene	13.87	228	145935	11.03	ng	99
82) Chrysene	13.91	228	90906	7.14	ng	98
85) Indeno(1,2,3-cd)pyrene	16.61	276	37655m	3.64	ng	
87) Benzo(b)fluoranthene	14.89	252	105199m	9.39	ng	
88) Benzo(k)fluoranthene	14.90	252	40779m	4.11	ng	
89) Benzo(a)pyrene	15.22	252	82157	8.28	ng	98
91) Benzo(g,h,i)perylene	17.01	276	38599	4.99	ng	97

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

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