

Data Path : Z:\HPCHEM1\BNA F\DATA\BF061316\
 Data File : BF087997.D
 Acq On : 14 Jun 2016 9:15
 Operator : UM/SJ
 Sample : H3395-01DL 2X
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 STA-928-PLATFORM(0-4)DL

Manual Integrations
 APPROVED

umangi
 6/14/2016 7:48:59 PM

Quant Time: Jun 14 19:23:40 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 13 14:10:49 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	125922	20.00	ng	0.01
21) Naphthalene-d8	8.08	136	482138	20.00	ng	0.00
38) Acenaphthene-d10	9.84	164	233284	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	350038	20.00	ng	0.01
75) Chrysene-d12	13.95	240	237842	20.00	ng	0.00
86) Perylene-d12	15.37	264	238385	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	459722	62.41	ng	0.00
7) Phenol-d6	6.43	99	636803	71.34	ng	0.00
23) Nitrobenzene-d5	7.36	82	345632	41.86	ng	0.00
41) 2,4,6-Tribromophenol	10.63	330	118780	48.22	ng	0.00
44) 2-Fluorobiphenyl	9.16	172	645493	40.57	ng	0.00
78) Terphenyl-d14	12.90	244	368020	35.21	ng	0.00
Target Compounds						
48) Acenaphthylene	9.70	152	171836	7.16	ng	Qvalue 97
49) Dimethylphthalate	9.55	163	45254	2.68	ng	99
70) Phenanthrene	11.35	178	229967	12.52	ng	97
71) Anthracene	11.39	178	138454	7.47	ng	98
74) Fluoranthene	12.54	202	942938	48.64	ng	97
77) Pyrene	12.77	202	823340	46.65	ng	100
80) Benzo(a)anthracene	13.94	228	535970	39.54	ng	# 89
82) Chrysene	13.98	228	357132m	28.01	ng	
85) Indeno(1,2,3-cd)pyrene	16.73	276	212074	17.90	ng	# 100
87) Benzo(b)fluoranthene	14.97	252	579582m	34.88	ng	
88) Benzo(k)fluoranthene	14.98	252	242151m	19.32	ng	
89) Benzo(a)pyrene	15.31	252	360810	26.84	ng	97
90) Dibenzo(a,h)anthracene	16.73	278	51896m	4.16	ng	
91) Benzo(a,h,i)perylene	17.14	276	172531	13.43	ng	97

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

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