

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060916\
 Data File : BF087853.D
 Acq On : 9 Jun 2016 15:33
 Operator : UM/SJ
 Sample : H3399-11 5X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SS-31

Quant Time: Jun 09 16:17:32 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 07 18:51:55 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	121011	20.00	ng	-0.02
21) Naphthalene-d8	8.10	136	447657	20.00	ng	-0.02
38) Acenaphthene-d10	9.85	164	215678	20.00	ng	-0.03
63) Phenanthrene-d10	11.32	188	353484	20.00	ng	-0.03
75) Chrysene-d12	13.95	240	223811	20.00	ng	-0.03
86) Perylene-d12	15.36	264	253179	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	225624	31.87	ng	-0.02
7) Phenol-d6	6.44	99	313704	36.57	ng	-0.03
23) Nitrobenzene-d5	7.38	82	167972	21.91	ng	-0.03
41) 2,4,6-Tribromophenol	10.64	330	56999	25.03	ng	-0.03
44) 2-Fluorobiphenyl	9.18	172	359876	22.12	ng	-0.03
78) Terphenyl-d14	12.91	244	223293	22.70	ng	-0.03
Target Compounds						Qvalue
70) Phenanthrene	11.36	178	334054	18.01	ng	97
71) Anthracene	11.40	178	91574	4.89	ng	97
74) Fluoranthene	12.54	202	372355	19.02	ng	98
77) Pyrene	12.76	202	304732	18.35	ng	98
80) Benzo(a)anthracene	13.94	228	149002	11.68	ng	100
82) Chrysene	13.98	228	99622	8.30	ng	96
85) Indeno(1,2,3-cd)pyrene	16.71	276	65060	5.84	ng	# 100
87) Benzo(b)fluoranthene	14.96	252	146294m	8.29	ng	
88) Benzo(k)fluoranthene	14.97	252	65944m	4.95	ng	
89) Benzo(a)pyrene	15.30	252	114483	8.02	ng	96
91) Benzo(g,h,i)perylene	17.12	276	59904	4.39	ng	97

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

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