

Data Path : Z:\HPCHEM1\BNA F\DATA\BF061016\
 Data File : BF087878.D
 Acq On : 10 Jun 2016 16:56
 Operator : UM/SJ
 Sample : H3426-11 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DUP-4

Manual Integrations
 APPROVED

SOHIL
 6/13/2016 7:13:19 PM

Quant Time: Jun 10 18:32:45 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	6.81	152	132423m	20.00	ng	-0.03
21) Naphthalene-d8	8.10	136	532620	20.00	ng	-0.02
38) Acenaphthene-d10	9.85	164	230458	20.00	ng	-0.03
63) Phenanthrene-d10	11.32	188	314579	20.00	ng	-0.03
75) Chrysene-d12	13.95	240	291693	20.00	ng	-0.03
86) Perylene-d12	15.36	264	244736	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	261645	33.78	ng	-0.03
7) Phenol-d6	6.44	99	360380	38.39	ng	-0.03
23) Nitrobenzene-d5	7.37	82	208228	22.83	ng	-0.05
41) 2,4,6-Tribromophenol	10.63	330	64186	26.38	ng	-0.05
44) 2-Fluorobiphenyl	9.18	172	416759	24.46	ng	-0.03
78) Terphenyl-d14	12.90	244	214009	16.70	ng	-0.05
Target Compounds						Qvalue
49) Dimethylphthalate	9.56	163	40157	2.41	ng	100
70) Phenanthrene	11.35	178	81082	4.91	ng	97
74) Fluoranthene	12.52	202	94443	5.42	ng	95
77) Pyrene	12.75	202	93095	4.30	ng	99
80) Benzo(a)anthracene	13.94	228	53574	3.22	ng	97
82) Chrysene	13.98	228	50877m	3.25	ng	
87) Benzo(b)fluoranthene	14.96	252	68762m	4.03	ng	
89) Benzo(a)pyrene	15.30	252	44211	3.20	ng	95

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

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