

Data Path : Z:\HPCHEM1\BNA F\DATA\BF062216\
 Data File : BF088266.D
 Acq On : 22 Jun 2016 23:12
 Operator : UM/SJ
 Sample : H3660-05
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-3(0-6)

Manual Integrations
 APPROVED

umangi
 6/23/2016 7:03:33 PM

Quant Time: Jun 23 11:31:17 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 22 14:55:14 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.60	152	81231	20.00	ng	0.00
21) Naphthalene-d8	7.88	136	335891	20.00	ng	-0.01
38) Acenaphthene-d10	9.63	164	146953	20.00	ng	-0.01
63) Phenanthrene-d10	11.12	188	227310	20.00	ng	0.00
75) Chrysene-d12	13.75	240	169632	20.00	ng	0.00
86) Perylene-d12	15.11	264	160274	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.19	112	510532	107.44	ng	0.01
7) Phenol-d6	6.26	99	609430	105.83	ng	0.00
23) Nitrobenzene-d5	7.18	82	384562	66.86	ng	0.00
41) 2,4,6-Tribromophenol	10.42	330	119764	77.18	ng	-0.01
44) 2-Fluorobiphenyl	8.97	172	681562	72.01	ng	0.00
78) Terphenyl-d14	12.70	244	448892	60.22	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.37	163	137355	12.90	ng	99
74) Fluoranthene	12.32	202	50028	3.97	ng	96
77) Pyrene	12.55	202	48017	3.81	ng	98
80) Benzo(a)anthracene	13.74	228	22179m	2.29	ng	
82) Chrysene	13.77	228	18904	2.08	ng	99
87) Benzo(b)fluoranthene	14.74	252	34488m	3.09	ng	
89) Benzo(a)pyrene	15.06	252	21774	2.41	ng	# 92

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

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