

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102516\
 Data File : BF090831.D
 Acq On : 26 Oct 2016 2:05
 Operator : UM/SJ
 Sample : H5396-04 2X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P2-S3

Manual Integrations
 APPROVED

umangi
 10/26/2016 4:46:14 PM

Quant Time: Oct 26 03:25:37 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 25 18:56:36 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	213388	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	896139	20.00	ng	0.00
38) Acenaphthene-d10	9.84	164	396077	20.00	ng	-0.01
63) Phenanthrene-d10	11.32	188	585855	20.00	ng	0.00
75) Chrysene-d12	13.96	240	401517	20.00	ng	0.00
86) Perylene-d12	15.38	264	370493	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	664938	50.25	ng	0.00
7) Phenol-d6	6.44	99	895386	49.96	ng	0.00
23) Nitrobenzene-d5	7.37	82	528545	32.35	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	153336	48.91	ng	-0.01
44) 2-Fluorobiphenyl	9.16	172	855461	33.53	ng	-0.01
78) Terphenyl-d14	12.91	244	602947	34.42	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.56	163	269183	10.82	ng	# 97
74) Fluoranthene	12.53	202	97665	3.38	ng	96
77) Pyrene	12.76	202	88270	3.15	ng	96
80) Benzo(a)anthracene	13.95	228	48857m	2.22	ng	
82) Chrysene	13.99	228	64548m	3.01	ng	
87) Benzo(b)fluoranthene	14.98	252	80963m	3.55	ng	

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

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