

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080516\
 Data File : BF089632.D
 Acq On : 5 Aug 2016 16:02
 Operator : UM/SJ
 Sample : H4292-05
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SS8-0-2IN

Manual Integrations
 APPROVED

umangi
 8/8/2016 9:12:38 AM

Quant Time: Aug 08 01:54:14 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF080416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 05 01:39:23 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.44	152	49626	20.00	ng	-0.01
21) Naphthalene-d8	9.46	136	210980	20.00	ng	-0.01
38) Acenaphthene-d10	12.29	164	95231	20.00	ng	-0.01
63) Phenanthrene-d10	14.69	188	161175	20.00	ng	-0.01
75) Chrysene-d12	18.38	240	101193	20.00	ng	-0.01
86) Perylene-d12	20.05	264	99356	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	412290	139.25	ng	0.04
7) Phenol-d6	6.99	99	539991	134.44	ng	0.01
23) Nitrobenzene-d5	8.35	82	347316	91.06	ng	-0.01
41) 2,4,6-Tribromophenol	13.59	330	102433	130.34	ng	0.00
44) 2-Fluorobiphenyl	11.26	172	594947	81.19	ng	0.00
78) Terphenyl-d14	17.05	244	369928	72.17	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	11.94	163	155035	20.77	ng	100
70) Phenanthrene	14.72	178	18325	2.09	ng	99
74) Fluoranthene	16.49	202	28552	3.17	ng	97
77) Pyrene	16.79	202	23656	2.64	ng	# 94
87) Benzo(b)fluoranthene	19.65	252	16557m	2.71	ng	
89) Benzo(a)pyrene	19.99	252	12583	2.20	ng	# 88

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

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