

Data Path : Z:\HPCHEM1\BNA F\DATA\BF123115\
 Data File : BF084219.D
 Acq On : 1 Jan 2016 2:30
 Operator : UM/SJ
 Sample : G4982-03
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S8-0530

Manual Integrations
 APPROVED

UMANGI
 1/4/2016 4:29:48 PM

Quant Time: Jan 02 02:34:56 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 31 14:15:57 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	283436	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	1109110	20.00	ng	0.00
38) Acenaphthene-d10	9.95	164	494010	20.00	ng	0.00
63) Phenanthrene-d10	11.42	188	778847	20.00	ng	0.00
75) Chrysene-d12	14.07	240	569150	20.00	ng	0.00
86) Perylene-d12	15.51	264	464403	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	1139328	65.02	ng	0.01
7) Phenol-d6	6.56	99	978834	44.48	ng	0.02
23) Nitrobenzene-d5	7.48	82	1817624	91.21	ng	0.01
41) 2,4,6-Tribromophenol	10.74	330	615370	123.38	ng	0.00
44) 2-Fluorobiphenyl	9.28	172	3030828	109.77	ng	0.01
78) Terphenyl-d14	13.01	244	1439736	56.47	ng	0.00
Target Compounds						
9) Benzaldehyde	6.45	77	673542	47.00	ng	Qvalue 97
22) Acetophenone	7.34	105	10159614	380.94	ng	# 86
28) bis(2-Chloroethoxy)methane	7.96	93	96632	4.15	ng	# 63
31) Naphthalene	8.21	128	422886	8.40	ng	100
32) Benzoic acid	7.95	122	82281	12.22	ng	# 71
37) 2-Methylnaphthalene	8.90	142	168035	4.65	ng	97
49) Dimethylphthalate	9.66	163	158646m	4.49	ng	

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

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