

Data Path : Z:\HPCHEM1\BNA F\DATA\BF013118\
 Data File : BF102670.D
 Acq On : 31 Jan 2018 22:19
 Operator : SJ/JU
 Sample : J1378-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RO-72

Manual Integrations
 APPROVED

Sohil
 2/1/2018 6:00:30 PM

Quant Time: Feb 01 12:20:44 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 31 18:38:00 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.71	152	105726	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	420118	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	164970	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	281836	20.00	ng	0.00
75) Chrysene-d12	13.93	240	1080322	20.00	ng	0.06
86) Perylene-d12	15.27	264	15192m	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.34	112	677114	97.11	ng	0.02
7) Phenol-d6	6.39	99	638868	76.00	ng	0.03
23) Nitrobenzene-d5	7.28	82	427600	58.49	ng	0.00
41) 2,4,6-Tribromophenol	10.55	330	138371	84.65	ng	0.01
44) 2-Fluorobiphenyl	9.07	172	747169	66.59	ng	0.00
78) Terphenyl-d14	12.83	244	605019	12.63	ng	0.02
Target Compounds						
9) Benzaldehyde	6.26	77	213611	50.265	ng	Qvalue 94
10) Phenol	6.40	94	500128	54.495	ng	76
11) bis(2-Chloroethyl)ether	6.45	93	122504m	17.551	ng	
15) Benzyl Alcohol	6.88	79	78456m	13.228	ng	
20) 3+4-Methylphenols	7.17	107	145290	19.460	ng	88
22) Acetophenone	7.13	105	34158	3.411	ng	# 95
31) Naphthalene	8.02	128	56403	2.689	ng	# 96
32) Benzoic acid	7.93	122	211871m	44.229	ng	
37) 2-Methylnaphthalene	8.70	142	47092	3.371	ng	93
49) Dimethylphthalate	9.47	163	110950	8.137	ng	# 87
55) 4-Nitrophenol	9.89	139	22345	9.733	ng	90
73) Di-n-butylphthalate	11.79	149	59936	2.932	ng	# 91

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