

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012818\
 Data File : BF102566.D
 Acq On : 28 Jan 2018 19:16
 Operator : SJ/JU
 Sample : J1290-05
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-48-9.5-10.0

Manual Integrations
 APPROVED

Sohil
 1/29/2018 5:48:05 PM

Quant Time: Jan 29 04:27:46 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jan 27 20:48:40 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.71	152	158523	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	649023	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	275233	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	404327	20.00	ng	0.00
75) Chrysene-d12	13.88	240	261555	20.00	ng	0.00
86) Perylene-d12	15.32	264	253402	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	1126347	107.73	ng	0.00
7) Phenol-d6	6.36	99	1345101	106.72	ng	0.00
23) Nitrobenzene-d5	7.28	82	850129	75.27	ng	0.00
41) 2,4,6-Tribromophenol	10.54	330	228402	83.75	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	1103923	58.97	ng	0.00
78) Terphenyl-d14	12.82	244	711618	61.33	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.37	94	55519	4.035	ng	98
31) Naphthalene	8.02	128	121401	3.746	ng	99
49) Dimethylphthalate	9.46	163	145389	6.391	ng	100
51) Acenaphthene	9.78	154	230072	13.219	ng	98
54) Dibenzofuran	9.95	168	147044	6.242	ng	99
57) Fluorene	10.29	166	241702	12.949	ng	98
70) Phenanthrene	11.26	178	914760	38.825	ng	99
71) Anthracene	11.31	178	308824	12.842	ng	99
72) Carbazole	11.47	167	87349	3.723	ng	98
74) Fluoranthene	12.46	202	1189467	49.506	ng	98
77) Pyrene	12.68	202	1010996	49.994	ng	99
80) Benzo(a)anthracene	13.87	228	427334m	26.537	ng	
82) Chrysene	13.90	228	397046m	24.280	ng	
85) Indeno(1,2,3-cd)pyrene	16.72	276	141592	10.484	ng	96
87) Benzo(b)fluoranthene	14.90	252	442105m	26.918	ng	
88) Benzo(k)fluoranthene	14.93	252	129472m	8.344	ng	
89) Benzo(a)pyrene	15.26	252	305428	20.619	ng	# 95
90) Dibenzo(a,h)anthracene	16.72	278	36268	3.023	ng	# 74
91) Benzo(a,h,i)perylene	17.14	276	135251	11.416	ng	# 91

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