

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012918\
 Data File : BF102589.D
 Acq On : 29 Jan 2018 11:26
 Operator : SJ/JU
 Sample : J1257-17
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-20(5-15)

Manual Integrations
 APPROVED

Sohil
 1/30/2018 3:36:21 PM

Quant Time: Jan 30 00:35:41 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	156754	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	623872	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	268115	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	355936	20.00	ng	0.00
75) Chrysene-d12	13.89	240	222078	20.00	ng	0.01
86) Perylene-d12	15.34	264	261393	20.00	ng	0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	1167071	112.89	ng	0.01
7) Phenol-d6	6.36	99	1344066	107.84	ng	0.00
23) Nitrobenzene-d5	7.29	82	823266	75.83	ng	0.00
41) 2,4,6-Tribromophenol	10.55	330	218044	82.08	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	1260368	69.12	ng	0.00
78) Terphenyl-d14	12.83	244	699769	71.03	ng	0.01
Target Compounds						Qvalue
10) Phenol	6.37	94	58342	4.288	ng	86
22) Acetophenone	7.10	105	37690	2.534	ng	# 59
31) Naphthalene	8.02	128	77269	2.481	ng	# 93
37) 2-Methylnaphthalene	8.71	142	117629	5.670	ng	97
49) Dimethylphthalate	9.47	163	200949	9.068	ng	100
57) Fluorene	10.30	166	37129	2.042	ng	# 93
70) Phenanthrene	11.26	178	322646	15.556	ng	99
71) Anthracene	11.32	178	74096m	3.500	ng	
74) Fluoranthene	12.46	202	334943	15.836	ng	98
77) Pyrene	12.69	202	309089	18.001	ng	99
80) Benzo(a)anthracene	13.89	228	141014	10.314	ng	# 94
82) Chrysene	13.92	228	138278m	9.959	ng	
83) Bis(2-ethylhexyl)phthalate	13.86	149	122626	10.679	ng	95
84) Di-n-octyl phthalate	14.47	149	57175	3.062	ng	# 74
85) Indeno(1,2,3-cd)pyrene	16.75	276	96242	8.393	ng	95
87) Benzo(b)fluoranthene	14.92	252	178145m	10.515	ng	
88) Benzo(k)fluoranthene	14.94	252	68290m	4.266	ng	
89) Benzo(a)pyrene	15.28	252	140629	9.203	ng	# 64
91) Benzo(g,h,i)perylene	17.17	276	88728	7.260	ng	# 82

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