

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012818\
 Data File : BF102569.D
 Acq On : 28 Jan 2018 20:37
 Operator : SJ/JU
 Sample : J1253-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-02(0-5)

Manual Integrations
 APPROVED

Sohil
 1/29/2018 5:48:28 PM

Quant Time: Jan 29 04:42:47 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	166765	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	662709	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	267477	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	361332	20.00	ng	0.00
75) Chrysene-d12	13.88	240	257120	20.00	ng	0.00
86) Perylene-d12	15.32	264	232625	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	1139334	103.59	ng	0.01
7) Phenol-d6	6.37	99	1384876	104.44	ng	0.00
23) Nitrobenzene-d5	7.29	82	854712	74.12	ng	0.00
41) 2,4,6-Tribromophenol	10.55	330	205843	77.67	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	1247377	68.57	ng	0.00
78) Terphenyl-d14	12.82	244	724221	63.50	ng	0.00
Target Compounds						
10) Phenol	6.37	94	58201	4.021	ng	Qvalue # 76
48) Acenaphthylene	9.61	152	109941	3.796	ng	98
49) Dimethylphthalate	9.47	163	194108	8.780	ng	99
57) Fluorene	10.29	166	68822	3.794	ng	98
70) Phenanthrene	11.26	178	770325	36.585	ng	99
71) Anthracene	11.31	178	202316	9.414	ng	99
72) Carbazole	11.47	167	57954	2.764	ng	97
74) Fluoranthene	12.46	202	1103717	51.404	ng	97
77) Pyrene	12.68	202	938548	47.212	ng	100
80) Benzo(a)anthracene	13.87	228	550545	34.778	ng	97
82) Chrysene	13.91	228	504511	31.384	ng	97
83) Bis(2-ethylhexyl)phthalate	13.85	149	65965	4.962	ng	99
85) Indeno(1,2,3-cd)pyrene	16.72	276	184350	13.886	ng	97
87) Benzo(b)fluoranthene	14.91	252	587948m	38.995	ng	
88) Benzo(k)fluoranthene	14.93	252	207006m	14.532	ng	
89) Benzo(a)pyrene	15.26	252	378554	27.838	ng	# 95
90) Dibenzo(a,h)anthracene	16.72	278	51721	4.695	ng	# 87
91) Benzo(g,h,i)perylene	17.14	276	164450	15.120	ng	93

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