

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF012918\
 Data File : BF102595.D
 Acq On : 29 Jan 2018 14:09
 Operator : SJ/JU
 Sample : J1253-15 2X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-08(5-15)

Manual Integrations
 APPROVED

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

Quant Time: Jan 30 01:02:18 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	6.72	152	176965	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	701468	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	281630	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	400147	20.00	ng	0.00
75) Chrysene-d12	13.89	240	294823	20.00	ng	0.01
86) Perylene-d12	15.33	264	237065	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	636831	54.56	ng	0.00
7) Phenol-d6	6.36	99	760323	54.04	ng	0.00
23) Nitrobenzene-d5	7.28	82	426042	34.90	ng	0.00
41) 2,4,6-Tribromophenol	10.55	330	105845	37.93	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	710725	37.11	ng	0.00
78) Terphenyl-d14	12.82	244	465169	35.57	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.37	94	34870	2.270	ng	91
31) Naphthalene	8.02	128	112911	3.224	ng	100
37) 2-Methylnaphthalene	8.71	142	52370	2.245	ng	99
49) Dimethylphthalate	9.47	163	139368	5.987	ng	99
51) Acenaphthene	9.79	154	122254	6.864	ng	99
54) Dibenzofuran	9.96	168	109009	4.523	ng	98
57) Fluorene	10.30	166	168446	8.819	ng	97
70) Phenanthrene	11.27	178	1476390	63.317	ng	99
71) Anthracene	11.32	178	419027	17.606	ng	100
72) Carbazole	11.47	167	165286	7.119	ng	98
74) Fluoranthene	12.46	202	1456411	61.250	ng	99
77) Pyrene	12.69	202	1478327	64.854	ng	100
80) Benzo(a)anthracene	13.88	228	674237	37.145	ng	99
82) Chrysene	13.92	228	642073m	34.834	ng	
85) Indeno(1,2,3-cd)pyrene	16.74	276	206618	13.573	ng	97
87) Benzo(b)fluoranthene	14.92	252	589258m	38.350	ng	
88) Benzo(k)fluoranthene	14.94	252	191066m	13.162	ng	
89) Benzo(a)pyrene	15.27	252	427972	30.883	ng	# 95
90) Dibenzo(a,h)anthracene	16.74	278	55743	4.966	ng	# 66
91) Benzo(g,h,i)perylene	17.17	276	212028	19.129	ng	95

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