

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122717\
 Data File : BF101781.D
 Acq On : 27 Dec 2017 23:42
 Operator : SJ/JU
 Sample : I7064-09 2X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SAMPLE-5

Manual Integrations
 APPROVED

Sohil
 12/28/2017 2:03:11 PM

Quant Time: Dec 28 05:56:48 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 26 10:48:01 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	153174	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	593527	20.00	ng	0.00
38) Acenaphthene-d10	9.83	164	251672	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	397578	20.00	ng	0.00
75) Chrysene-d12	13.97	240	283204	20.00	ng	0.00
86) Perylene-d12	15.45	264	284983	20.00	ng	0.02

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.42	112	576911	56.10	ng	0.01
7) Phenol-d6	6.44	99	711771	55.62	ng	0.00
23) Nitrobenzene-d5	7.36	82	418309	39.23	ng	0.00
41) 2,4,6-Tribromophenol	10.63	330	124352	51.28	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	737292	45.44	ng	0.00
78) Terphenyl-d14	12.90	244	505478	43.03	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
10) Phenol	6.46	94	32790	2.248	ng	# 66
31) Naphthalene	8.09	128	268387	8.982	ng	99
37) 2-Methylnaphthalene	8.79	142	161302	8.286	ng	98
45) 1,1'-Biphenyl	9.25	154	61895	2.773	ng	98
48) Acenaphthylene	9.69	152	61604	2.284	ng	# 97
49) Dimethylphthalate	9.55	163	63913	3.157	ng	99
51) Acenaphthene	9.86	154	178449	11.011	ng	100
54) Dibenzofuran	10.03	168	411797	19.995	ng	99
57) Fluorene	10.38	166	271135	15.562	ng	99
70) Phenanthrene	11.36	178	2061932	93.258	ng	98
71) Anthracene	11.40	178	821978	36.395	ng	100
72) Carbazole	11.56	167	254118	12.018	ng	100
74) Fluoranthene	12.55	202	1695452	73.056	ng	98
77) Pyrene	12.77	202	1478104	69.543	ng	100
79) Butylbenzylphthalate	13.37	149	22535	2.343	ng	96
80) Benzo(a)anthracene	13.97	228	798867	42.719	ng	100
82) Chrysene	14.00	228	675429	38.254	ng	98
83) Bis(2-ethylhexyl)phthalate	13.92	149	46247	3.797	ng	99
85) Indeno(1,2,3-cd)pyrene	16.94	276	263090	16.733	ng	# 89
87) Benzo(b)fluoranthene	15.02	252	730712m	42.067	ng	
88) Benzo(k)fluoranthene	15.04	252	233536m	13.828	ng	
89) Benzo(a)pyrene	15.39	252	526827	32.900	ng	96
90) Dibenzo(a,h)anthracene	16.93	278	60690m	4.414	ng	
91) Benzo(g,h,i)perylene	17.39	276	216106	15.874	ng	95

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

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