

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122217\
 Data File : BF101659.D
 Acq On : 22 Dec 2017 17:53
 Operator : SJ/JU
 Sample : I6683-11 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-02-122017-C

Manual Integrations
 APPROVED

Sohil
 12/26/2017 1:37:22 PM

Quant Time: Dec 22 23:26:44 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 22 16:07:58 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	160025	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	605506	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	253516	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	428326	20.00	ng	0.00
75) Chrysene-d12	13.97	240	299874	20.00	ng	0.00
86) Perylene-d12	15.43	264	312268	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	505408	47.05	ng	0.00
7) Phenol-d6	6.43	99	619783	46.36	ng	0.00
23) Nitrobenzene-d5	7.36	82	376164	34.58	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	110857	45.38	ng	0.00
44) 2-Fluorobiphenyl	9.14	172	593671	36.33	ng	0.00
78) Terphenyl-d14	12.90	244	445841	35.84	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	8.09	128	298109	9.779	ng	98
37) 2-Methylnaphthalene	8.78	142	193330	9.734	ng	98
48) Acenaphthylene	9.69	152	77632	2.857	ng	# 97
49) Dimethylphthalate	9.54	163	69499	3.408	ng	99
51) Acenaphthene	9.86	154	141554	8.671	ng	99
54) Dibenzofuran	10.03	168	178067	8.583	ng	# 95
57) Fluorene	10.37	166	284667	16.220	ng	99
70) Phenanthrene	11.35	178	1394913	58.561	ng	98
71) Anthracene	11.39	178	530177	21.790	ng	99
72) Carbazole	11.56	167	95822	4.206	ng	97
74) Fluoranthene	12.54	202	1500409	60.011	ng	98
77) Pyrene	12.77	202	1389898	61.758	ng	99
80) Benzo(a)anthracene	13.96	228	753055	38.031	ng	99
82) Chrysene	13.99	228	587602	31.429	ng	96
85) Indeno(1,2,3-cd)pyrene	16.90	276	263732	15.841	ng	# 91
87) Benzo(b)fluoranthene	15.01	252	648080m	34.050	ng	
88) Benzo(k)fluoranthene	15.03	252	248230m	13.414	ng	
89) Benzo(a)pyrene	15.37	252	515138	29.359	ng	98
90) Dibenzo(a,h)anthracene	16.90	278	67077	4.453	ng	# 89
91) Benzo(g,h,i)perylene	17.35	276	244098	16.363	ng	93

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

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