

Data Path : Z:\HPCHEM1\BNA F\DATA\BF111617\
 Data File : BF100629.D
 Acq On : 16 Nov 2017 14:37
 Operator : SJ/JU
 Sample : I6302-01 10X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HR-04-111317-A

Manual Integrations
 APPROVED

SOHIL
 11/17/2017 1:24:43 PM

Quant Time: Nov 16 15:08:37 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 16 11:34:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	106850	20.00	ng	-0.02
21) Naphthalene-d8	8.16	136	424747	20.00	ng	-0.02
38) Acenaphthene-d10	9.92	164	198019	20.00	ng	-0.02
63) Phenanthrene-d10	11.40	188	365257	20.00	ng	-0.02
75) Chrysene-d12	14.04	240	214264	20.00	ng	-0.02
86) Perylene-d12	15.52	264	250961	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	78358	11.76	ng	-0.02
7) Phenol-d6	6.49	99	97139	11.82	ng	-0.03
23) Nitrobenzene-d5	7.43	82	54464	8.48	ng	-0.04
41) 2,4,6-Tribromophenol	10.70	330	12532	6.21	ng	-0.02
44) 2-Fluorobiphenyl	9.23	172	125113	9.62	ng	-0.03
78) Terphenyl-d14	12.98	244	102456	10.99	ng	-0.03
Target Compounds						Qvalue
31) Naphthalene	8.19	128	309455	15.126	ng	99
37) 2-Methylnaphthalene	8.87	142	150075	11.049	ng	97
45) 1,1'-Biphenyl	9.33	154	43776	2.556	ng	95
48) Acenaphthylene	9.77	152	74510	3.619	ng	# 96
51) Acenaphthene	9.95	154	237321	18.951	ng	99
54) Dibenzofuran	10.12	168	274612	15.646	ng	95
57) Fluorene	10.46	166	405958	31.573	ng	99
70) Phenanthrene	11.44	178	1834553	95.394	ng	99
71) Anthracene	11.48	178	719973	36.901	ng	99
72) Carbazole	11.63	167	162048	9.001	ng	98
74) Fluoranthene	12.63	202	1696817	83.253	ng	98
77) Pylene	12.86	202	1482569	97.068	ng	100
80) Benzo(a)anthracene	14.03	228	744489	57.699	ng	97
82) Chrysene	14.07	228	594460	47.577	ng	97
85) Indeno(1,2,3-cd)pyrene	17.00	276	361106m	35.731	ng	
87) Benzo(b)fluoranthene	15.09	252	737310m	49.590	ng	
88) Benzo(k)fluoranthene	15.12	252	303674m	21.617	ng	
89) Benzo(a)pyrene	15.46	252	589621	44.732	ng	96
90) Dibenzo(a,h)anthracene	17.01	278	95674m	8.875	ng	
91) Benzo(a,h,i)perylene	17.46	276	312946	29.657	ng	93

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

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