

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF032116\
 Data File : BF085657.D
 Acq On : 22 Mar 2016 9:53
 Operator : UM/SJ
 Sample : H1815-10DL 5X
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-07-COMPDL

Manual Integrations
 APPROVED

Sohil
 3/22/2016 5:34:48 PM

Quant Time: Mar 22 11:51:41 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF031816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 18 23:31:39 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	132875m	20.00	ng	-0.02
21) Naphthalene-d8	8.13	136	557037	20.00	ng	-0.01
38) Acenaphthene-d10	9.88	164	215564	20.00	ng	-0.02
63) Phenanthrene-d10	11.36	188	345687	20.00	ng	-0.02
75) Chrysene-d12	14.00	240	272332	20.00	ng	-0.01
86) Perylene-d12	15.42	264	285955	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.43	112	188791	23.86	ng	-0.02
7) Phenol-d6	6.47	99	228175	22.49	ng	-0.02
23) Nitrobenzene-d5	7.40	82	130642	13.61	ng	-0.03
41) 2,4,6-Tribromophenol	10.66	330	32932	15.60	ng	-0.02
44) 2-Fluorobiphenyl	9.20	172	202846	12.14	ng	-0.02
78) Terphenyl-d14	12.94	244	103159	9.12	ng	-0.02
Target Compounds						Qvalue
57) Fluorene	10.42	166	41694	2.98	ng	98
70) Phenanthrene	11.39	178	505199	27.72	ng	99
71) Anthracene	11.43	178	99902	5.23	ng	99
72) Carbazole	11.59	167	43357	2.63	ng	100
74) Fluoranthene	12.57	202	576009	30.18	ng	94
77) Pyrene	12.80	202	527856	26.99	ng	100
80) Benzo(a)anthracene	13.99	228	253856	16.06	ng	94
82) Chrysene	14.03	228	235955	15.51	ng	97
85) Indeno(1,2,3-cd)pyrene	16.81	276	107981	8.50	ng	94
87) Benzo(b)fluoranthene	15.02	252	297103m	15.92	ng	
88) Benzo(k)fluoranthene	15.04	252	94206m	6.13	ng	
89) Benzo(a)pyrene	15.36	252	214921	13.59	ng	98
90) Dibenzo(a,h)anthracene	16.81	278	32813	2.49	ng	# 76
91) Benzo(g,h,i)perylene	17.23	276	100361	7.86	ng	98

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

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