

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021116\
 Data File : BF085018.D
 Acq On : 11 Feb 2016 14:53
 Operator : SJ/IZ
 Sample : H1361-02DL 10X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-1DL

Manual Integrations
 APPROVED

Sohil
 2/12/2016 12:26:09 PM

Quant Time: Feb 11 17:55:15 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 10 13:28:51 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.60	152	127372	20.00	ng	-0.02
21) Naphthalene-d8	7.90	136	557728	20.00	ng	-0.02
38) Acenaphthene-d10	9.64	164	261927	20.00	ng	-0.02
63) Phenanthrene-d10	11.12	188	441740	20.00	ng	-0.02
75) Chrysene-d12	13.76	240	294242	20.00	ng	-0.01
86) Perylene-d12	15.12	264	280134	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.19	112	139781	17.09	ng	-0.01
7) Phenol-d6	6.26	99	183773	18.13	ng	-0.02
23) Nitrobenzene-d5	7.18	82	121400	12.26	ng	-0.02
41) 2,4,6-Tribromophenol	10.43	330	42030	15.38	ng	-0.02
44) 2-Fluorobiphenyl	8.97	172	191330	8.57	ng	-0.02
78) Terphenyl-d14	12.71	244	151774	11.69	ng	-0.02
Target Compounds						Qvalue
51) Acenaphthene	9.68	154	51990	3.06	ng	93
57) Fluorene	10.19	166	43029	2.48	ng	97
70) Phenanthrene	11.15	178	563212	22.99	ng	99
71) Anthracene	11.20	178	276722	11.21	ng	98
72) Carbazole	11.36	167	70780	3.29	ng	97
74) Fluoranthene	12.33	202	1334479	53.81	ng	98
77) Pyrene	12.56	202	1258355	55.86	ng	99
80) Benzo(a)anthracene	13.75	228	619487m	33.92	ng	
82) Chrysene	13.78	228	527087m	29.76	ng	
85) Indeno(1,2,3-cd)pyrene	16.37	276	226967	16.28	ng	95
87) Benzo(b)fluoranthene	14.75	252	605202m	32.15	ng	
88) Benzo(k)fluoranthene	14.77	252	166118m	10.67	ng	
89) Benzo(a)pyrene	15.06	252	393956	24.56	ng	# 94
90) Dibenzo(a,h)anthracene	16.37	278	72299	5.36	ng	# 91
91) Benzo(g,h,i)perylene	16.73	276	203537	14.97	ng	98

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

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