

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122316\
 Data File : BF091930.D
 Acq On : 23 Dec 2016 18:11
 Operator : UM/SJ
 Sample : H6238-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PIT-002-COMP

Quant Time: Dec 26 10:20:21 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 21 16:17:51 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	205879	20.00	ng	-0.01
21) Naphthalene-d8	8.03	136	770227	20.00	ng	-0.01
38) Acenaphthene-d10	9.79	164	359258	20.00	ng	0.00
63) Phenanthrene-d10	11.28	188	526446	20.00	ng	0.01
75) Chrysene-d12	13.89	240	436754	20.00	ng	-0.01
86) Perylene-d12	15.29	264	422860	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	1463618	118.70	ng	0.03
7) Phenol-d6	6.43	99	1615524	107.32	ng	0.02
23) Nitrobenzene-d5	7.31	82	1069619	77.22	ng	-0.01
41) 2,4,6-Tribromophenol	10.59	330	314824	105.89	ng	0.01
44) 2-Fluorobiphenyl	9.12	172	1393107	78.46	ng	0.00
78) Terphenyl-d14	12.84	244	959113	53.26	ng	-0.01
Target Compounds						Qvalue
10) Phenol	6.44	94	86784	5.06	ng	88
49) Dimethylphthalate	9.52	163	335599	14.61	ng	# 96
51) Acenaphthene	9.82	154	112954	5.56	ng	# 62
57) Fluorene	10.34	166	224986	11.27	ng	# 85
70) Phenanthrene	11.30	178	646709	22.66	ng	# 93
74) Fluoranthene	12.48	202	147587m	2.26	ng	
77) Pyrene	12.69	202	235676m	7.35	ng	
80) Benzo(a)anthracene	13.88	228	61013	2.47	ng	# 80
82) Chrysene	13.92	228	62244	2.52	ng	# 90
85) Indeno(1,2,3-cd)pyrene	16.61	276	61365	2.86	ng	96
87) Benzo(b)fluoranthene	14.89	252	133574m	5.07	ng	
89) Benzo(a)pyrene	15.22	252	76968	3.29	ng	# 87
91) Benzo(a,h,i)perylene	17.01	276	57151	2.86	ng	95

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

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