

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011316\
 Data File : BF084481.D
 Acq On : 14 Jan 2016 10:35
 Operator : SJ/UM
 Sample : H1077-05 5X
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 POST-EX-5-20160108

Quant Time: Jan 14 16:57:27 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 14 14:15:52 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	377886	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	1498381	20.00	ng	0.00
38) Acenaphthene-d10	9.87	164	597709	20.00	ng	0.00
63) Phenanthrene-d10	11.36	188	924499	20.00	ng	0.00
75) Chrysene-d12	14.00	240	720477	20.00	ng	0.00
86) Perylene-d12	15.41	264	696702	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	578816	24.78	ng	0.00
7) Phenol-d6	6.46	99	685526	23.36	ng	-0.01
23) Nitrobenzene-d5	7.39	82	400673	14.88	ng	-0.01
41) 2,4,6-Tribromophenol	10.66	330	116645	19.33	ng	-0.01
44) 2-Fluorobiphenyl	9.20	172	763913	17.46	ng	0.00
78) Terphenyl-d14	12.93	244	414180	12.83	ng	-0.01
Target Compounds						Qvalue
31) Naphthalene	8.14	128	430966	6.34	ng	98
37) 2-Methylnaphthalene	8.83	142	109433	2.24	ng	97
48) Acenaphthylene	9.73	152	139407	2.53	ng	98
49) Dimethylphthalate	9.60	163	118962	2.78	ng	# 91
51) Acenaphthene	9.90	154	160078	4.33	ng	97
57) Fluorene	10.42	166	226217	3.94	ng	98
70) Phenanthrene	11.38	178	1624129	33.59	ng	97
71) Anthracene	11.42	178	432428	9.12	ng	99
72) Carbazole	11.58	167	167979	3.62	ng	97
74) Fluoranthene	12.57	202	1677467	33.21	ng	98
77) Pyrene	12.80	202	1526728	27.00	ng	98
80) Benzo(a)anthracene	13.99	228	801033	18.93	ng	94
82) Chrysene	14.02	228	745096m	18.33	ng	
85) Indeno(1,2,3-cd)pyrene	16.80	276	344649	10.70	ng	95
87) Benzo(b)fluoranthene	15.01	252	925856m	20.32	ng	
88) Benzo(k)fluoranthene	15.04	252	271549m	7.29	ng	
89) Benzo(a)pyrene	15.36	252	651314	16.85	ng	# 95
90) Dibenzo(a,h)anthracene	16.80	278	79966	2.40	ng	# 90
91) Benzo(a,h,i)perylene	17.21	276	326491	9.48	ng	97

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

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