

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121616\
 Data File : BF091687.D
 Acq On : 16 Dec 2016 21:54
 Operator : UM/SJ
 Sample : H6099-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SS03-0-2IN

Quant Time: Dec 19 12:57:34 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 16 00:35:41 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	273473	20.00	ng	-0.01
21) Naphthalene-d8	8.05	136	1062559	20.00	ng	-0.02
38) Acenaphthene-d10	9.81	164	563663	20.00	ng	-0.01
63) Phenanthrene-d10	11.29	188	789750	20.00	ng	-0.02
75) Chrysene-d12	13.92	240	425127	20.00	ng	-0.02
86) Perylene-d12	15.32	264	507899	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	1575932	97.09	ng	0.00
7) Phenol-d6	6.42	99	1398024	69.08	ng	-0.01
23) Nitrobenzene-d5	7.33	82	1517711	79.72	ng	-0.02
41) 2,4,6-Tribromophenol	10.60	330	419535	89.61	ng	-0.01
44) 2-Fluorobiphenyl	9.14	172	2079718	63.75	ng	-0.01
78) Terphenyl-d14	12.88	244	1321174	70.52	ng	-0.02
Target Compounds						Qvalue
10) Phenol	6.43	94	101027	4.27	ng	88
32) Benzoic acid	7.78	122	16145	6.19	ng	92
49) Dimethylphthalate	9.53	163	304006	8.52	ng	# 98
70) Phenanthrene	11.31	178	336145	8.60	ng	100
74) Fluoranthene	12.50	202	451020	11.01	ng	98
77) Pyrene	12.73	202	331416	9.83	ng	98
80) Benzo(a)anthracene	13.91	228	142211	5.76	ng	98
82) Chrysene	13.94	228	100679m	3.90	ng	
83) Bis(2-ethylhexyl)phthalate	13.92	149	33846	2.01	ng	# 92
85) Indeno(1,2,3-cd)pyrene	16.66	276	80412	3.46	ng	96
87) Benzo(b)fluoranthene	14.92	252	158555m	5.24	ng	
89) Benzo(a)pyrene	15.27	252	131746m	4.83	ng	
91) Benzo(a,h,i)perylene	17.06	276	77117	3.10	ng	# 96

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

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