

Data Path : Z:\HPCHEM1\BNA F\DATA\BF071516\
 Data File : BF089017.D
 Acq On : 15 Jul 2016 15:25
 Operator : UM/SJ
 Sample : H3972-09
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LSS-SB-SB10(0-2ft)

Quant Time: Jul 15 23:23:27 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 15 23:12:46 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.67	152	65762	20.00	ng	0.00
21) Naphthalene-d8	7.96	136	257629	20.00	ng	0.00
38) Acenaphthene-d10	9.71	164	99814	20.00	ng	0.00
63) Phenanthrene-d10	11.19	188	150738	20.00	ng	0.00
75) Chrysene-d12	13.81	240	114303	20.00	ng	-0.01
86) Perylene-d12	15.18	264	82470	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.28	112	538258	122.67	ng	0.01
7) Phenol-d6	6.34	99	839871	148.13	ng	0.01
23) Nitrobenzene-d5	7.24	82	420094	85.45	ng	0.00
41) 2,4,6-Tribromophenol	10.50	330	116436	125.62	ng	0.00
44) 2-Fluorobiphenyl	9.04	172	723616	114.51	ng	0.00
78) Terphenyl-d14	12.77	244	412056	80.29	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.44	163	177265	24.93	ng	99
70) Phenanthrene	11.21	178	72628	8.81	ng	97
71) Anthracene	11.26	178	25812	3.15	ng	99
74) Fluoranthene	12.39	202	130804	15.66	ng	97
77) Pyrene	12.62	202	126827	13.09	ng	97
80) Benzo(a)anthracene	13.79	228	69708	9.47	ng	92
82) Chrysene	13.83	228	49141m	6.75	ng	
85) Indeno(1,2,3-cd)pyrene	16.45	276	22219	3.97	ng	# 100
87) Benzo(b)fluoranthene	14.80	252	61371m	11.86	ng	
88) Benzo(k)fluoranthene	14.82	252	20084m	4.46	ng	
89) Benzo(a)pyrene	15.12	252	42483	9.70	ng	98
90) Dibenzo(a,h)anthracene	16.45	278	9652	2.64	ng	# 82
91) Benzo(a,h,i)perylene	16.82	276	23209	6.13	ng	98

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

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