

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060716\
 Data File : BF087802.D
 Acq On : 7 Jun 2016 23:19
 Operator : UM/SJ
 Sample : H3379-07
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SS-4

Manual Integrations
 APPROVED

sohil
 6/8/2016 7:13:23 PM

Quant Time: Jun 08 05:15:37 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 07 18:51:55 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	99512	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	400991	20.00	ng	0.00
38) Acenaphthene-d10	9.87	164	174763	20.00	ng	-0.01
63) Phenanthrene-d10	11.36	188	284286	20.00	ng	0.00
75) Chrysene-d12	13.98	240	199723	20.00	ng	-0.01
86) Perylene-d12	15.41	264	209524	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	705395	121.18	ng	0.01
7) Phenol-d6	6.48	99	954144	135.25	ng	0.00
23) Nitrobenzene-d5	7.40	82	550743	80.20	ng	-0.01
41) 2,4,6-Tribromophenol	10.66	330	207729	112.57	ng	-0.01
44) 2-Fluorobiphenyl	9.20	172	999738	90.20	ng	-0.01
78) Terphenyl-d14	12.94	244	679497	77.42	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.60	163	253083	19.99	ng	98
51) Acenaphthene	9.91	154	25878	2.31	ng	98
70) Phenanthrene	11.38	178	368253	24.69	ng	98
71) Anthracene	11.43	178	100096	6.65	ng	98
72) Carbazole	11.58	167	31164	2.21	ng	98
74) Fluoranthene	12.56	202	461879	29.34	ng	100
77) Pyrene	12.79	202	446084	30.10	ng	99
80) Benzo(a)anthracene	13.98	228	233469	20.51	ng	99
82) Chrysene	14.01	228	174022m	16.25	ng	
85) Indeno(1,2,3-cd)pyrene	16.77	276	106916	10.75	ng	# 100
87) Benzo(b)fluoranthene	14.99	252	238437m	16.33	ng	
88) Benzo(k)fluoranthene	15.01	252	99398m	9.02	ng	
89) Benzo(a)pyrene	15.34	252	183590	15.54	ng	99
90) Dibenzo(a,h)anthracene	16.79	278	26666	2.43	ng	92
91) Benzo(a,h,i)perylene	17.19	276	98998	8.77	ng	98

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

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