

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051616\
 Data File : BF087092.D
 Acq On : 16 May 2016 20:22
 Operator : UM/SJ
 Sample : H3038-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WEST-ABOUTMENT-N.SIDE

Manual Integrations
 APPROVED

umangi
 5/17/2016 7:16:05 PM

Quant Time: May 17 13:31:05 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 17 01:12:04 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.60	152	151534	20.00	ng	0.00
21) Naphthalene-d8	7.90	136	582795	20.00	ng	0.00
38) Acenaphthene-d10	9.64	164	237014	20.00	ng	0.00
63) Phenanthrene-d10	11.12	188	387529	20.00	ng	0.00
75) Chrysene-d12	13.75	240	309068	20.00	ng	-0.01
86) Perylene-d12	15.11	264	245705	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.20	112	908134	101.49	ng	0.00
7) Phenol-d6	6.27	99	1159798	96.98	ng	-0.01
23) Nitrobenzene-d5	7.18	82	746493	64.32	ng	-0.01
41) 2,4,6-Tribromophenol	10.43	330	212296	84.61	ng	-0.01
44) 2-Fluorobiphenyl	8.97	172	1099325	77.95	ng	-0.01
78) Terphenyl-d14	12.71	244	920938	63.22	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.37	163	128806	7.12	ng	Qvalue 100
70) Phenanthrene	11.14	178	429942	20.79	ng	99
71) Anthracene	11.20	178	100782	4.76	ng	98
74) Fluoranthene	12.33	202	646843	30.41	ng	92
77) Pyrene	12.55	202	561959	23.75	ng	97
80) Benzo(a)anthracene	13.74	228	306722	17.68	ng	99
82) Chrysene	13.77	228	226053	12.81	ng	99
85) Indeno(1,2,3-cd)pyrene	16.35	276	94711	6.45	ng	# 91
87) Benzo(b)fluoranthene	14.74	252	259577m	16.26	ng	
88) Benzo(k)fluoranthene	14.75	252	104884m	8.02	ng	
89) Benzo(a)pyrene	15.06	252	182445	13.25	ng	97
90) Dibenzo(a,h)anthracene	16.37	278	31945	2.75	ng	# 90
91) Benzo(a,h,i)perylene	16.73	276	83219	7.06	ng	96

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

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