

Data Path : Z:\HPCHEM1\BNA F\DATA\BF070316\
 Data File : BF088685.D
 Acq On : 4 Jul 2016 10:11
 Operator : UM/SJ
 Sample : H3769-01
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EPCH-B1(00-0.5)

Manual Integrations

umangi
 7/5/2016 7:50:42 PM

Quant Time: Jul 05 17:07:24 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 05 01:06:52 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	117418	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	403643	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	168719	20.00	ng	0.00
63) Phenanthrene-d10	11.30	188	271146	20.00	ng	0.01
75) Chrysene-d12	13.93	240	196225	20.00	ng	0.01
86) Perylene-d12	15.33	264	172997	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	606557	77.42	ng	0.01
7) Phenol-d6	6.42	99	747903	73.88	ng	0.00
23) Nitrobenzene-d5	7.35	82	477870	62.04	ng	0.00
41) 2,4,6-Tribromophenol	10.60	330	131885	84.18	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	861397	80.65	ng	0.00
78) Terphenyl-d14	12.88	244	495888	56.29	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.54	163	102752	8.55	ng	98
70) Phenanthrene	11.32	178	327267	22.08	ng	98
71) Anthracene	11.37	178	89842	6.10	ng	98
72) Carbazole	11.52	167	34235	2.47	ng	98
74) Fluoranthene	12.50	202	494215	32.89	ng	97
77) Pyrene	12.73	202	477001	28.67	ng	99
80) Benzo(a)anthracene	13.92	228	269701	21.35	ng	97
82) Chrysene	13.95	228	222888m	17.83	ng	
85) Indeno(1,2,3-cd)pyrene	16.65	276	106661	11.09	ng	# 100
87) Benzo(b)fluoranthene	14.93	252	308264m	28.41	ng	
88) Benzo(k)fluoranthene	14.95	252	45836m	4.85	ng	
89) Benzo(a)pyrene	15.27	252	182331m	19.84	ng	
90) Dibenzo(a,h)anthracene	16.66	278	27019	3.52	ng	# 90
91) Benzo(a,h,i)perylene	17.05	276	97837	12.32	ng	98

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

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