

Data Path : Z:\HPCHEM1\BNA F\DATA\BF070316\
 Data File : BF088687.D
 Acq On : 4 Jul 2016 11:09
 Operator : UM/SJ
 Sample : H3769-06
 Misc :
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EPCH-B3(00-0.5)

Manual Integrations

umangi
 7/5/2016 7:50:36 PM

Quant Time: Jul 05 17:15:09 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	104426	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	359338	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	143449	20.00	ng	0.00
63) Phenanthrene-d10	11.30	188	236255	20.00	ng	0.01
75) Chrysene-d12	13.92	240	184354	20.00	ng	0.00
86) Perylene-d12	15.33	264	143651	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	597139	85.70	ng	0.01
7) Phenol-d6	6.42	99	724102	80.43	ng	0.00
23) Nitrobenzene-d5	7.35	82	443791	64.72	ng	0.00
41) 2,4,6-Tribromophenol	10.60	330	117008	87.84	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	808668	89.05	ng	0.00
78) Terphenyl-d14	12.88	244	496375	59.97	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.54	163	96031	9.40	ng	98
70) Phenanthrene	11.32	178	73280	5.67	ng	97
74) Fluoranthene	12.50	202	193593	14.79	ng	98
77) Pyrene	12.73	202	190403	12.18	ng	98
80) Benzo(a)anthracene	13.92	228	107928	9.09	ng	96
82) Chrysene	13.95	228	81961m	6.98	ng	
83) Bis(2-ethylhexyl)phthalate	13.92	149	15980	2.01	ng	# 95
85) Indeno(1,2,3-cd)pyrene	16.66	276	33322	3.69	ng	# 100
87) Benzo(b)fluoranthene	14.93	252	94492m	10.49	ng	
88) Benzo(k)fluoranthene	14.95	252	20275m	2.58	ng	
89) Benzo(a)pyrene	15.27	252	63454	8.32	ng	95
91) Benzo(a,h,i)perylene	17.06	276	33156	5.03	ng	# 92

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

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