

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF070216\
 Data File : BF088635.D
 Acq On : 2 Jul 2016 23:58
 Operator : UM/SJ
 Sample : H3769-01
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

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

Manual Integrations

umangi
 7/5/2016 7:35:47 PM

Quant Time: Jul 03 05:00:12 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF063016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 30 18:01:55 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	113873	20.00	ng	-0.01
21) Naphthalene-d8	8.08	136	464737	20.00	ng	-0.02
38) Acenaphthene-d10	9.83	164	186951	20.00	ng	-0.02
63) Phenanthrene-d10	11.31	188	307774	20.00	ng	-0.01
75) Chrysene-d12	13.94	240	225957	20.00	ng	-0.01
86) Perylene-d12	15.34	264	196437	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	725972	95.55	ng	-0.01
7) Phenol-d6	6.43	99	960715	97.86	ng	-0.01
23) Nitrobenzene-d5	7.36	82	587220	66.22	ng	-0.02
41) 2,4,6-Tribromophenol	10.62	330	153396	88.36	ng	-0.02
44) 2-Fluorobiphenyl	9.15	172	854702	72.22	ng	-0.02
78) Terphenyl-d14	12.89	244	542011	53.43	ng	-0.02
Target Compounds						Qvalue
49) Dimethylphthalate	9.55	163	109417	8.22	ng	99
70) Phenanthrene	11.34	178	376694	22.39	ng	99
71) Anthracene	11.38	178	96514	5.77	ng	97
72) Carbazole	11.53	167	38197	2.43	ng	98
74) Fluoranthene	12.51	202	552686	32.40	ng	96
77) Pyrene	12.74	202	502672	26.24	ng	99
80) Benzo(a)anthracene	13.93	228	292292	20.09	ng	97
82) Chrysene	13.97	228	256326m	17.81	ng	
85) Indeno(1,2,3-cd)pyrene	16.67	276	122605	11.07	ng	# 100
87) Benzo(b)fluoranthene	14.95	252	321809m	26.12	ng	
88) Benzo(k)fluoranthene	14.96	252	83028m	7.74	ng	
89) Benzo(a)pyrene	15.28	252	203746	19.53	ng	98
90) Dibenzo(a,h)anthracene	16.69	278	31891	3.66	ng	# 86
91) Benzo(g,h,i)perylene	17.09	276	113551	12.60	ng	97

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

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