

Data Path : Z:\HPCHEM1\BNA F\DATA\BF070516\
 Data File : BF088729.D
 Acq On : 6 Jul 2016 4:01
 Operator : UM/SJ
 Sample : H3881-16
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 Q6-B5-0-11-C

Manual Integrations

sohil
 7/6/2016 7:20:32 PM

Quant Time: Jul 06 13:02:06 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 06 01:47:10 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.78	152	102680	20.00	ng	0.00
21) Naphthalene-d8	8.06	136	361232	20.00	ng	-0.01
38) Acenaphthene-d10	9.80	164	156086	20.00	ng	-0.01
63) Phenanthrene-d10	11.28	188	252295	20.00	ng	-0.01
75) Chrysene-d12	13.91	240	204268	20.00	ng	-0.01
86) Perylene-d12	15.30	264	182042	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	575325	83.97	ng	0.00
7) Phenol-d6	6.41	99	652167	73.67	ng	0.00
23) Nitrobenzene-d5	7.34	82	404941	58.75	ng	-0.01
41) 2,4,6-Tribromophenol	10.59	330	127235	87.78	ng	-0.01
44) 2-Fluorobiphenyl	9.14	172	730596	73.94	ng	0.00
78) Terphenyl-d14	12.87	244	504982	55.06	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.53	163	140132	12.60	ng	99
70) Phenanthrene	11.30	178	112394	8.15	ng	99
71) Anthracene	11.36	178	30302	2.21	ng	98
74) Fluoranthene	12.49	202	165754	11.86	ng	95
77) Pyrene	12.72	202	134279	7.75	ng	97
80) Benzo(a)anthracene	13.90	228	69370m	5.27	ng	
82) Chrysene	13.93	228	70894m	5.45	ng	
85) Indeno(1,2,3-cd)pyrene	16.62	276	33508	3.35	ng	# 100
87) Benzo(b)fluoranthene	14.90	252	97314m	8.52	ng	
88) Benzo(k)fluoranthene	14.93	252	20436m	2.06	ng	
89) Benzo(a)pyrene	15.23	252	58109	6.01	ng	98
91) Benzo(a,h,i)perylene	17.02	276	34783	4.16	ng	98

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

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