

Data Path : Z:\HPCHEM1\BNA F\DATA\BF070516\
 Data File : BF088728.D
 Acq On : 6 Jul 2016 3:32
 Operator : UM/SJ
 Sample : H3881-15
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 Q6-B8-0-5-C

Manual Integrations

sohil
 7/6/2016 7:20:29 PM

Quant Time: Jul 06 12:59:00 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	101130	20.00	ng	0.00
21) Naphthalene-d8	8.06	136	363767	20.00	ng	-0.01
38) Acenaphthene-d10	9.80	164	159276	20.00	ng	-0.01
63) Phenanthrene-d10	11.28	188	250921	20.00	ng	-0.01
75) Chrysene-d12	13.91	240	195738	20.00	ng	-0.01
86) Perylene-d12	15.30	264	177571	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	574442	85.13	ng	0.00
7) Phenol-d6	6.41	99	693889	79.58	ng	0.00
23) Nitrobenzene-d5	7.34	82	436452	62.88	ng	-0.01
41) 2,4,6-Tribromophenol	10.59	330	135095	91.34	ng	-0.01
44) 2-Fluorobiphenyl	9.14	172	769773	76.34	ng	0.00
78) Terphenyl-d14	12.87	244	485496	55.25	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.53	163	151628	13.36	ng	99
70) Phenanthrene	11.30	178	89349	6.51	ng	98
74) Fluoranthene	12.49	202	167018	12.01	ng	95
77) Pyrene	12.72	202	129186	7.78	ng	97
80) Benzo(a)anthracene	13.90	228	69041m	5.48	ng	
82) Chrysene	13.93	228	71311m	5.72	ng	
85) Indeno(1,2,3-cd)pyrene	16.62	276	30871	3.22	ng	# 100
87) Benzo(b)fluoranthene	14.91	252	98244m	8.82	ng	
88) Benzo(k)fluoranthene	14.93	252	20920m	2.16	ng	
89) Benzo(a)pyrene	15.23	252	51066	5.41	ng	96
91) Benzo(g,h,i)perylene	17.02	276	31017	3.81	ng	98

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

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