

Data Path : Z:\HPCHEM1\BNA F\DATA\BF070516\
 Data File : BF088727.D
 Acq On : 6 Jul 2016 3:02
 Operator : UM/SJ
 Sample : H3881-13
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 Q6-B6-0-5-C

Manual Integrations

sohil
 7/6/2016 7:20:28 PM

Quant Time: Jul 06 12:57:00 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	107785	20.00	ng	0.00
21) Naphthalene-d8	8.06	136	388869	20.00	ng	-0.01
38) Acenaphthene-d10	9.80	164	168505	20.00	ng	-0.01
63) Phenanthrene-d10	11.28	188	281773	20.00	ng	-0.01
75) Chrysene-d12	13.91	240	206894	20.00	ng	-0.01
86) Perylene-d12	15.30	264	196098	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	605917	84.25	ng	0.00
7) Phenol-d6	6.41	99	852022	91.69	ng	0.00
23) Nitrobenzene-d5	7.34	82	516284	69.58	ng	-0.01
41) 2,4,6-Tribromophenol	10.59	330	154530	98.76	ng	-0.01
44) 2-Fluorobiphenyl	9.14	172	920698	86.31	ng	0.00
78) Terphenyl-d14	12.87	244	561151	60.41	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.53	163	182285	15.18	ng	99
70) Phenanthrene	11.30	178	58325	3.79	ng	99
74) Fluoranthene	12.50	202	126571	8.11	ng	93
77) Pyrene	12.72	202	122113	6.96	ng	97
80) Benzo(a)anthracene	13.90	228	67721m	5.08	ng	
82) Chrysene	13.93	228	78598m	5.96	ng	
83) Bis(2-ethylhexyl)phthalate	13.91	149	54428	6.09	ng	99
85) Indeno(1,2,3-cd)pyrene	16.63	276	35804	3.53	ng	# 100
87) Benzo(b)fluoranthene	14.91	252	105923m	8.61	ng	
88) Benzo(k)fluoranthene	14.93	252	24583m	2.30	ng	
89) Benzo(a)pyrene	15.25	252	56879	5.46	ng	97
91) Benzo(a,h,i)perylene	17.02	276	36588	4.07	ng	99

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

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