

Data Path : Z:\HPCHEM1\BNA F\DATA\BF070516\
 Data File : BF088726.D
 Acq On : 6 Jul 2016 2:33
 Operator : UM/SJ
 Sample : H3881-11
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 Q5-B1-0-11-C

Manual Integrations

sohil
 7/6/2016 7:20:26 PM

Quant Time: Jul 06 12:54:57 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	107537	20.00	ng	0.00
21) Naphthalene-d8	8.06	136	383722	20.00	ng	-0.01
38) Acenaphthene-d10	9.80	164	168854	20.00	ng	-0.01
63) Phenanthrene-d10	11.28	188	264613	20.00	ng	-0.01
75) Chrysene-d12	13.91	240	197864	20.00	ng	-0.01
86) Perylene-d12	15.30	264	193099	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	619720	86.37	ng	0.00
7) Phenol-d6	6.41	99	847682	91.43	ng	0.00
23) Nitrobenzene-d5	7.34	82	513348	70.11	ng	-0.01
41) 2,4,6-Tribromophenol	10.59	330	159286	101.59	ng	-0.01
44) 2-Fluorobiphenyl	9.14	172	875464	81.90	ng	0.00
78) Terphenyl-d14	12.87	244	553746	62.33	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.53	163	183376	15.24	ng	98
70) Phenanthrene	11.30	178	142980	9.88	ng	98
71) Anthracene	11.36	178	36667	2.55	ng	97
74) Fluoranthene	12.49	202	131088	8.94	ng	95
77) Pyrene	12.71	202	116007	6.91	ng	96
80) Benzo(a)anthracene	13.90	228	58243	4.57	ng	99
82) Chrysene	13.93	228	37555	2.98	ng	96
85) Indeno(1,2,3-cd)pyrene	16.62	276	22098	2.28	ng	# 100
87) Benzo(b)fluoranthene	14.90	252	63813m	5.27	ng	
89) Benzo(a)pyrene	15.23	252	39050	3.81	ng	96
91) Benzo(g,h,i)perylene	17.02	276	21966	2.48	ng	98

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

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