

Data Path : Z:\HPCHEM1\BNA F\DATA\BF072116\
 Data File : BF089174.D
 Acq On : 21 Jul 2016 12:14
 Operator : UM/SJ
 Sample : H4112-04
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 KSB-03-0.2-2.0

Manual Integrations

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Quant Time: Jul 22 01:18:09 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 20 19:03:15 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.60	152	53324	20.00	ng	0.00
21) Naphthalene-d8	7.88	136	224518	20.00	ng	0.00
38) Acenaphthene-d10	9.62	164	92275	20.00	ng	-0.01
63) Phenanthrene-d10	11.10	188	163483	20.00	ng	-0.01
75) Chrysene-d12	13.71	240	104569	20.00	ng	-0.01
86) Perylene-d12	15.06	264	96866	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.19	112	246791	70.39	ng	0.00
7) Phenol-d6	6.25	99	335525	74.21	ng	-0.02
23) Nitrobenzene-d5	7.16	82	210858	49.90	ng	-0.01
41) 2,4,6-Tribromophenol	10.41	330	43566	52.44	ng	-0.01
44) 2-Fluorobiphenyl	8.96	172	369756	55.21	ng	-0.01
78) Terphenyl-d14	12.69	244	253937	52.67	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.36	163	79751	11.26	ng	98
74) Fluoranthene	12.31	202	36502	3.87	ng	93
77) Pyrene	12.53	202	32881	3.81	ng	99
80) Benzo(a)anthracene	13.71	228	18682m	2.81	ng	
82) Chrysene	13.75	228	14896m	2.34	ng	
85) Indeno(1,2,3-cd)pyrene	16.29	276	12073	2.61	ng	94
87) Benzo(b)fluoranthene	14.71	252	24596m	3.51	ng	
89) Benzo(a)pyrene	15.02	252	18577m	3.41	ng	
91) Benzo(g,h,i)perylene	16.65	276	10993	2.54	ng	# 90

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

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