

Data Path : Z:\HPCHEM1\BNA F\DATA\BF110716\
 Data File : BF091073.D
 Acq On : 7 Nov 2016 23:07
 Operator : UM/SJ
 Sample : H5526-15DL 5X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-105-0.5-1.0DL

Manual Integrations
 APPROVED

sohil
 11/8/2016 4:32:32 PM

Quant Time: Nov 08 01:23:00 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 08 00:06:22 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.67	152	130358	20.00	ng	0.00
21) Naphthalene-d8	7.96	136	441989	20.00	ng	0.00
38) Acenaphthene-d10	9.71	164	225828	20.00	ng	0.00
63) Phenanthrene-d10	11.20	188	353467	20.00	ng	0.00
75) Chrysene-d12	13.83	240	371870	20.00	ng	0.00
86) Perylene-d12	15.22	264	336075	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.25	112	209092	26.49	ng	0.00
7) Phenol-d6	6.33	99	283894	26.50	ng	0.00
23) Nitrobenzene-d5	7.24	82	67042	8.27	ng	0.00
41) 2,4,6-Tribromophenol	10.50	330	37028	18.14	ng	-0.01
44) 2-Fluorobiphenyl	9.04	172	255961	19.51	ng	-0.01
78) Terphenyl-d14	12.77	244	233653	15.68	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.44	163	73937	4.99	ng	99
70) Phenanthrene	11.22	178	111026	5.91	ng	97
74) Fluoranthene	12.41	202	551973	28.32	ng	88
77) Pyrene	12.63	202	587958	24.14	ng	96
80) Benzo(a)anthracene	13.83	228	440003	20.84	ng	98
82) Chrysene	13.86	228	377099m	18.11	ng	
85) Indeno(1,2,3-cd)pyrene	16.51	276	327435	18.96	ng	# 91
87) Benzo(b)fluoranthene	14.83	252	688333m	30.20	ng	
88) Benzo(k)fluoranthene	14.84	252	296819m	14.84	ng	
89) Benzo(a)pyrene	15.16	252	486252	24.50	ng	97
90) Dibenzo(a,h)anthracene	16.52	278	88249m	5.14	ng	
91) Benzo(a,h,i)perylene	16.91	276	284760	18.36	ng	# 90

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

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