

Data Path : Z:\HPCHEM1\BNA F\DATA\BF070316\
 Data File : BF088690.D
 Acq On : 4 Jul 2016 12:36
 Operator : UM/SJ
 Sample : H3822-03
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 D10-B3(6-12)C

Manual Integrations

umangi
 7/5/2016 7:50:02 PM

Quant Time: Jul 05 01:36:45 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 05 01:06:52 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	110501	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	388686	20.00	ng	0.00
38) Acenaphthene-d10	9.82	164	162987	20.00	ng	0.00
63) Phenanthrene-d10	11.30	188	266243	20.00	ng	0.01
75) Chrysene-d12	13.92	240	195882	20.00	ng	0.00
86) Perylene-d12	15.33	264	168175	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	918549	124.58	ng	0.02
7) Phenol-d6	6.43	99	1213926	127.42	ng	0.01
23) Nitrobenzene-d5	7.36	82	741820	100.02	ng	0.01
41) 2,4,6-Tribromophenol	10.60	330	167917	110.95	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	1084086	105.06	ng	0.00
78) Terphenyl-d14	12.88	244	595364	67.70	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.54	163	197564	17.01	ng	99
70) Phenanthrene	11.32	178	65889	4.53	ng	97
74) Fluoranthene	12.50	202	129863	8.80	ng	99
77) Pyrene	12.73	202	152043	9.15	ng	98
80) Benzo(a)anthracene	13.91	228	71345	5.66	ng	97
82) Chrysene	13.95	228	58639	4.70	ng	98
85) Indeno(1,2,3-cd)pyrene	16.65	276	22599	2.35	ng	# 100
87) Benzo(b)fluoranthene	14.93	252	60120m	5.70	ng	
88) Benzo(k)fluoranthene	14.94	252	23885m	2.60	ng	
89) Benzo(a)pyrene	15.26	252	43594	4.88	ng	98
91) Benzo(g,h,i)perylene	17.05	276	22877	2.96	ng	96

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

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