

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121815\
 Data File : BF083900.D
 Acq On : 18 Dec 2015 6:40
 Operator : UM/SJ
 Sample : G4759-01
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 K210-0-2

Manual Integrations
 APPROVED

MMdadoda
 12/18/2015 6:22:02 PM

Quant Time: Dec 18 07:42:13 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 18 01:01:02 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	73793	20.00	ng	0.00
21) Naphthalene-d8	8.16	136	261730	20.00	ng	0.00
38) Acenaphthene-d10	9.92	164	119530	20.00	ng	0.00
63) Phenanthrene-d10	11.41	188	147638	20.00	ng	0.02
75) Chrysene-d12	14.08	240	111180	20.00	ng	0.05
86) Perylene-d12	15.54	264	95576	20.00	ng	0.07
System Monitoring Compounds						
5) 2-Fluorophenol	5.49	112	356903	75.69	ng	0.01
7) Phenol-d6	6.52	99	427230	74.02	ng	0.01
23) Nitrobenzene-d5	7.44	82	240758	52.50	ng	0.00
41) 2,4,6-Tribromophenol	10.72	330	86209	65.47	ng	0.01
44) 2-Fluorobiphenyl	9.24	172	519627	58.05	ng	0.00
78) Terphenyl-d14	13.01	244	251041	50.23	ng	0.03
Target Compounds						
49) Dimethylphthalate	9.63	163	126223	12.58	ng	99
70) Phenanthrene	11.44	178	24858	2.79	ng	# 69
74) Fluoranthene	12.65	202	33652	3.78	ng	63
77) Pyrene	12.88	202	49320	3.96	ng	# 89
80) Benzo(a)anthracene	14.07	228	32177	4.71	ng	# 73
82) Chrysene	14.10	228	24949m	3.78	ng	
85) Indeno(1,2,3-cd)pyrene	16.99	276	29207	5.91	ng	97
87) Benzo(b)fluoranthene	15.12	252	44637m	6.41	ng	
88) Benzo(k)fluoranthene	15.13	252	14616m	2.54	ng	
91) Benzo(g,h,i)perylene	17.43	276	56176	12.80	ng	# 89

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

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