

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042115\
 Data File : BF078700.D
 Acq On : 21 Apr 2015 23:42
 Operator : TP/IZ
 Sample : G1961-02
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SW-017

Manual Integrations
 APPROVED

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 4/22/2015 5:22:53 PM

Quant Time: Apr 22 03:52:33 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 22 00:52:13 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	5.82	152	54450	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	245305	20.00	ng	-0.01
38) Acenaphthene-d10	11.17	164	134773	20.00	ng	-0.01
63) Phenanthrene-d10	13.90	188	262287	20.00	ng	-0.01
75) Chrysene-d12	17.77	240	256421	20.00	ng	-0.01
86) Perylene-d12	19.45	264	235856	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	3.84	112	194916	59.02	ng	0.01
7) Phenol-d6	5.32	99	254129	61.40	ng	0.00
23) Nitrobenzene-d5	6.75	82	153224	37.86	ng	-0.01
41) 2,4,6-Tribromophenol	12.64	330	67226	60.14	ng	-0.01
44) 2-Fluorobiphenyl	9.99	172	343449	36.43	ng	-0.01
78) Terphenyl-d14	16.43	244	373494	32.91	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	10.80	163	21607	2.13	ng	# 99
70) Phenanthrene	13.94	178	33582	2.21	ng	98
74) Fluoranthene	15.83	202	88676	5.54	ng	95
77) Pyrene	16.14	202	80324	4.99	ng	97
80) Benzo(a)anthracene	17.76	228	54002	3.54	ng	98
82) Chrysene	17.79	228	53891	3.76	ng	98
85) Indeno(1,2,3-cd)pyrene	20.58	276	34885	2.14	ng	# 82
87) Benzo(b)fluoranthene	19.04	252	72625m	4.98	ng	
89) Benzo(a)pyrene	19.38	252	48744	3.83	ng	# 91
91) Benzo(g,h,i)perylene	20.90	276	32110	2.51	ng	96

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

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