

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022421\
 Data File : BN013751.D
 Acq On : 25 Feb 2021 06:32
 Operator : CG/JU
 Sample : M1473-03
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GW803-20-25-022321-FD

Quant Time: Feb 25 07:05:48 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN022421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 24 21:36:22 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.741	152	5629	0.40	ng	0.00
7) Naphthalene-d8	10.530	136	22969	0.40	ng	# 0.00
13) Acenaphthene-d10	14.372	164	13526	0.40	ng	-0.01
19) Phenanthrene-d10	17.128	188	29182	0.40	ng	# 0.00
29) Chrysene-d12	21.300	240	30437	0.40	ng	-0.01
36) Perylene-d12	23.555	264	30649	0.40	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.333	112	1668	0.11	ng	0.00
5) Phenol-d6	6.903	99	1392	0.07	ng	0.00
8) Nitrobenzene-d5	8.883	82	4271	0.29	ng	0.00
11) 2-Methylnaphthalene-d10	12.120	152	8294	0.22	ng	0.00
14) 2,4,6-Tribromophenol	15.869	330	2049	0.37	ng	0.00
15) 2-Fluorobiphenyl	13.001	172	12711	0.27	ng	-0.01
27) Fluoranthene-d10	19.156	212	25341	0.29	ng	0.00
31) Terphenyl-d14	19.762	244	26401	0.39	ng	0.00
Target Compounds						
						Qvalue
9) Naphthalene	10.581	128	2566	0.05	ng	# 86
12) 2-Methylnaphthalene	12.195	142	1311	0.03	ng	# 90
17) Acenaphthene	14.435	154	1216	0.03	ng	# 88
18) Fluorene	15.425	166	2394	0.05	ng	# 97
24) Pentachlorophenol	16.775	266	216	0.04	ng	99
25) Phenanthrene	17.165	178	25927	0.34	ng	99
26) Anthracene	17.250	178	3554	0.05	ng	# 95
28) Fluoranthene	19.186	202	8033	0.09	ng	# 97
30) Pyrene	19.548	202	8305	0.09	ng	96
32) Benzo(a)anthracene	21.290	228	4142	0.04	ng	92
33) Chrysene	21.343	228	4571	0.05	ng	# 96
34) Bis(2-ethylhexyl)phtha...	21.237	149	12107	0.27	ng	99
35) Indeno(1,2,3-cd)pyrene	25.824	276	3876	0.03	ng	99
37) Benzo(b)fluoranthene	22.877	252	4571	0.05	ng	# 79
38) Benzo(k)fluoranthene	22.922	252	4104	0.04	ng	# 39
39) Benzo(a)pyrene	23.452	252	2915	0.03	ng	# 68
40) Dibenzo(a,h)anthracene	25.842	278	2818	0.03	ng	# 62
41) Benzo(g,h,i)perylene	26.509	276	3377	0.04	ng	# 85

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

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