

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022421\
 Data File : BN013754.D
 Acq On : 25 Feb 2021 08:19
 Operator : CG/JU
 Sample : M1473-07
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GWBR1-119-129-022221

Quant Time: Feb 25 09:34:59 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	7.740	152	5322	0.40	ng	0.00
7) Naphthalene-d8	10.530	136	20889	0.40	ng	0.00
13) Acenaphthene-d10	14.371	164	11946	0.40	ng	-0.01
19) Phenanthrene-d10	17.128	188	26443	0.40	ng	# 0.00
29) Chrysene-d12	21.303	240	26907	0.40	ng	-0.01
36) Perylene-d12	23.551	264	26681	0.40	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.332	112	1876	0.13	ng	0.00
5) Phenol-d6	6.903	99	1464	0.08	ng	0.00
8) Nitrobenzene-d5	8.883	82	4221	0.31	ng	0.00
11) 2-Methylnaphthalene-d10	12.120	152	8229	0.24	ng	0.00
14) 2,4,6-Tribromophenol	15.869	330	2159	0.44	ng	0.00
15) 2-Fluorobiphenyl	13.001	172	12839	0.30	ng	-0.01
27) Fluoranthene-d10	19.153	212	24201	0.31	ng	0.00
31) Terphenyl-d14	19.759	244	56399	0.95	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.263	88	1811	0.28	ng	# 64
9) Naphthalene	10.568	128	2499	0.05	ng	# 87
12) 2-Methylnaphthalene	12.191	142	904	0.03	ng	# 84
25) Phenanthrene	17.165	178	3551	0.05	ng	100
26) Anthracene	17.250	178	1579	0.02	ng	# 96
28) Fluoranthene	19.183	202	3294	0.04	ng	# 93
30) Pyrene	19.546	202	3388	0.04	ng	97
32) Benzo(a)anthracene	21.292	228	2277	0.03	ng	# 89
33) Chrysene	21.335	228	2166	0.03	ng	# 90
34) Bis(2-ethylhexyl)phtha...	21.239	149	9750	0.25	ng	99
35) Indeno(1,2,3-cd)pyrene	25.820	276	2003	0.02	ng	95
37) Benzo(b)fluoranthene	22.876	252	2138	0.03	ng	# 65
38) Benzo(k)fluoranthene	22.921	252	1935	0.02	ng	# 27
39) Benzo(a)pyrene	23.451	252	1754	0.02	ng	# 63
40) Dibenzo(a,h)anthracene	25.837	278	1645	0.02	ng	# 47
41) Benzo(g,h,i)perylene	26.511	276	1688	0.02	ng	# 78

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

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