

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022521\
 Data File : BN013765.D
 Acq On : 25 Feb 2021 13:11
 Operator : CG/JU
 Sample : M1473-03
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

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

Quant Time: Feb 25 13:41:19 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN022421.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 25 11:31:11 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.741	152	5516	0.40	ng	0.00
7) Naphthalene-d8	10.530	136	22169	0.40	ng	0.00
13) Acenaphthene-d10	14.372	164	13057	0.40	ng	0.00
19) Phenanthrene-d10	17.128	188	27988	0.40	ng	# 0.00
29) Chrysene-d12	21.300	240	28901	0.40	ng	0.00
36) Perylene-d12	23.548	264	28714	0.40	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.333	112	1675	0.11	ng	0.00
5) Phenol-d6	6.903	99	1416	0.07	ng	0.00
8) Nitrobenzene-d5	8.883	82	4654	0.33	ng	0.00
11) 2-Methylnaphthalene-d10	12.120	152	8052	0.22	ng	0.00
14) 2,4,6-Tribromophenol	15.869	330	1868	0.35	ng	0.00
15) 2-Fluorobiphenyl	13.001	172	12346	0.27	ng	0.00
27) Fluoranthene-d10	19.156	212	24056	0.29	ng	0.00
31) Terphenyl-d14	19.757	244	24284	0.38	ng	0.00
Target Compounds						Qvalue
9) Naphthalene	10.568	128	2546	0.05	ng	# 86
12) 2-Methylnaphthalene	12.191	142	1283	0.03	ng	# 88
17) Acenaphthene	14.435	154	1207	0.03	ng	# 92
18) Fluorene	15.425	166	2618	0.06	ng	# 89
24) Pentachlorophenol	16.775	266	196	0.04	ng	97
25) Phenanthrene	17.164	178	25383	0.35	ng	99
26) Anthracene	17.250	178	3406	0.05	ng	# 96
28) Fluoranthene	19.186	202	7746	0.09	ng	# 96
30) Pyrene	19.543	202	7901	0.09	ng	97
32) Benzo(a)anthracene	21.290	228	3946	0.04	ng	# 90
33) Chrysene	21.343	228	4348	0.05	ng	# 92
34) Bis(2-ethylhexyl)phtha...	21.237	149	12490	0.30	ng	100
35) Indeno(1,2,3-cd)pyrene	25.814	276	3642	0.03	ng	97
37) Benzo(b)fluoranthene	22.874	252	4382	0.05	ng	# 72
38) Benzo(k)fluoranthene	22.922	252	3835	0.04	ng	# 49
39) Benzo(a)pyrene	23.449	252	2893	0.04	ng	# 59
40) Dibenzo(a,h)anthracene	25.835	278	2702	0.03	ng	# 62
41) Benzo(g,h,i)perylene	26.502	276	3191	0.04	ng	# 79

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

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