Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN120621\
Data File : BN017733.D
Acq On : 06 Dec 2021 16:17
Operator : CG/JU
Sample : PB141179BL
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
PB141179BL

Quant Time: Dec 06 16:56:55 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN120621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 06 15:58:05 2021
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.761	152	18370	0.400	ng	0.00
7) Naphthalene-d8	10.545	136	72937	0.400	ng	0.01
13) Acenaphthene-d10	14.373	164	44469	0.400	ng	0.00
19) Phenanthrene-d10	17.117	188	86808	0.400	ng	# 0.00
29) Chrysene-d12	21.314	240	57663	0.400	ng	# 0.01
35) Perylene-d12	23.620	264	40032	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.391	112	12483	0.317	ng	0.04
5) Phenol-d6	6.950	99	11033	0.261	ng	0.02
8) Nitrobenzene-d5	8.897	82	13639	0.277	ng	0.00
11) 2-Methylnaphthalene-d10	12.134	152	36596	0.274	ng	0.00
14) 2,4,6-Tribromophenol	15.883	330	1902	0.207	ng	0.01
15) 2-Fluorobiphenyl	13.003	172	46541	0.283	ng	0.00
27) Fluoranthene-d10	19.154	212	78628	0.269	ng	0.00
31) Terphenyl-d14	19.760	244	59977	0.405	ng	0.00
Target Compounds						Qvalue
6) bis(2-Chloroethyl)ether	7.189	93	80	0.002	ng	91
9) Naphthalene	10.596	128	294	0.002	ng	# 34
10) Hexachlorobutadiene	10.874	225	22	0.000	ng	# 95
12) 2-Methylnaphthalene	12.205	142	115	0.001	ng	# 73
16) Acenaphthylene	14.094	152	106	0.001	ng	# 79
17) Acenaphthene	14.436	154	69	0.001	ng	88
18) Fluorene	15.439	166	73	0.000	ng	# 97
21) 4-Bromophenyl-phenylether	16.326	248	16	0.000	ng	# 74
22) Hexachlorobenzene	16.435	284	29	0.000	ng	# 13
25) Phenanthrene	17.166	178	197	0.001	ng	# 77
26) Anthracene	17.263	178	59	0.000	ng	# 80
28) Fluoranthene	19.184	202	224	0.001	ng	# 65
30) Pyrene	19.546	202	140	0.001	ng	# 73
32) Benzo(a)anthracene	21.314	228	285	0.002	ng	# 28
33) Chrysene	21.314	228	211	0.001	ng	# 33
36) Indeno(1,2,3-cd)pyrene	26.006	276	41	0.000	ng	# 50
37) Benzo(b)fluoranthene	22.942	252	947	0.007	ng	# 1
38) Benzo(k)fluoranthene	22.942	252	714	0.006	ng	# 1
39) Benzo(a)pyrene	23.541	252	1029	0.009	ng	# 1
40) Dibenzo(a,h)anthracene	26.026	278	5	0.000	ng	# 1
41) Benzo(g,h,i)perylene	26.738	276	101	0.001	ng	# 1

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

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