Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010121\
Data File : BM028315.D
Acq On : 31 Dec 2020 21:03
Operator : JU/CG
Sample : PB133672BS
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : PB133672BS

Quant Time: Jan 01 00:00:57 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-SIM-BM010121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 31 16:18:03 2020
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units		Dev(Min)
Internal Standards							
1) 1,4-Dichlorobenzene-d4	8.135	152	725	0.40	ng		0.00
7) Naphthalene-d8	10.959	136	2880	0.40	ng		0.00
13) Acenaphthene-d10	14.763	164	1489	0.40	ng		0.00
19) Phenanthrene-d10	17.480	188	2940	0.40	ng	#	0.00
27) Chrysene-d12	21.600	240	2785	0.40	ng	#	-0.01
34) Perylene-d12	23.934	264	2603	0.40	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.638	112	867	0.35	ng		0.00
5) Phenol-d6	7.273	99	1165	0.35	ng		0.00
8) Nitrobenzene-d5	9.299	82	915	0.35	ng		0.00
11) 2-Methylnaphthalene-d10	12.531	152	1744	0.35	ng		0.00
14) 2,4,6-Tribromophenol	16.237	330	259	0.33	ng		0.00
15) 2-Fluorobiphenyl	13.390	172	2274	0.37	ng		0.00
25) Fluoranthene-d10	19.486	212	3024	0.33	ng		0.00
29) Terphenyl-d14	20.061	244	2587	0.38	ng		-0.01
Target Compounds							Qvalue
2) 1,4-Dioxane	3.407	88	359	0.35	ng		87
3) n-Nitrosodimethylamine	3.762	42	545	0.36	ng	#	94
6) bis(2-Chloroethyl)ether	7.547	93	851	0.36	ng		95
9) Naphthalene	11.010	128	2941	0.35	ng		98
10) Hexachlorobutadiene	11.289	225	585	0.36	ng	#	100
12) 2-Methylnaphthalene	12.606	142	1969	0.35	ng		96
16) Acenaphthylene	14.481	152	2300	0.35	ng		98
17) Acenaphthene	14.824	154	1686	0.35	ng		95
18) Fluorene	15.804	166	2114	0.35	ng		100
20) Hexachlorobenzene	16.800	284	714	0.36	ng		94
21) Atrazine	19.516	200	654	0.33	ng	#	36
22) Pentachlorophenol	17.140	266	265	0.32	ng		98
23) Phenanthrene	17.523	178	3197	0.35	ng		100
24) Anthracene	17.619	178	2696	0.34	ng		99
26) Fluoranthene	19.516	202	3352	0.33	ng	#	96
28) Pyrene	19.873	202	3360	0.35	ng		99
30) Benzo(a)anthracene	21.590	228	3156	0.34	ng		98
31) Chrysene	21.643	228	3419	0.36	ng		98
32) Bis(2-ethylhexyl)phtha...	21.494	149	1483	0.34	ng		95
33) Indeno(1,2,3-cd)pyrene	26.310	276	3716	0.36	ng	#	91
35) Benzo(b)fluoranthene	23.229	252	3639	0.38	ng	#	90
36) Benzo(k)fluoranthene	23.274	252	3282	0.36	ng	#	89
37) Benzo(a)pyrene	23.831	252	3091	0.37	ng	#	87
38) Dibenzo(a,h)anthracene	26.323	278	3237	0.37	ng	#	88
39) Benzo(g,h,i)perylene	27.042	276	3237	0.37	ng	#	86

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

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